

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102523\
 Data File : BF135940.D
 Acq On : 25 Oct 2023 13:35
 Operator : CG\JU
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICCC040

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 10/26/2023
 Supervised By :mohammad ahmed 10/26/2023

Quant Time: Oct 26 03:51:20 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 26 03:12:14 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.828	152	174615	20.000	ng	0.00
21) Naphthalene-d8	8.104	136	679462	20.000	ng	0.00
39) Acenaphthene-d10	9.863	164	343421	20.000	ng	0.00
64) Phenanthrene-d10	11.351	188	636222	20.000	ng	0.00
76) Chrysene-d12	14.010	240	405550	20.000	ng	0.00
86) Perylene-d12	15.492	264	388236	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.451	112	891092	78.802	ng	0.00
7) Phenol-d6	6.457	99	1116671	79.320	ng	0.00
23) Nitrobenzene-d5	7.392	82	824244	79.104	ng	0.00
42) 2,4,6-Tribromophenol	10.657	330	267231	88.901	ng	0.00
45) 2-Fluorobiphenyl	9.186	172	1709929	79.728	ng	0.00
79) Terphenyl-d14	12.957	244	2051702	84.526	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.616	88	209242	39.834	ng	100
3) Pyridine	3.369	79	578708	39.915	ng	100
4) n-Nitrosodimethylamine	3.340	42	251281	39.748	ng	100
6) Aniline	6.492	93	750491	39.591	ng	100
8) 2-Chlorophenol	6.610	128	466148	39.820	ng	100
9) Benzaldehyde	6.381	77	300536	39.464	ng	100
10) Phenol	6.469	94	575385m	39.093	ng	
11) bis(2-Chloroethyl)ether	6.569	93	450695	39.342	ng	100
12) 1,3-Dichlorobenzene	6.763	146	487526	39.208	ng	100
13) 1,4-Dichlorobenzene	6.845	146	493092	39.485	ng	100
14) 1,2-Dichlorobenzene	6.998	146	455152	39.249	ng	100
15) Benzyl Alcohol	6.969	79	403223	40.517	ng	100
16) 2,2'-oxybis(1-Chloropr...	7.104	45	662249	39.254	ng	100
17) 2-Methylphenol	7.075	107	408232	39.567	ng	100
18) Hexachloroethane	7.334	117	168415	39.661	ng	100
19) n-Nitroso-di-n-propyla...	7.239	70	317229	39.626	ng	100
20) 3+4-Methylphenols	7.228	107	496790	39.895	ng	100
22) Acetophenone	7.234	105	617027	40.207	ng	100
24) Nitrobenzene	7.410	77	449839	40.409	ng	100
25) Isophorone	7.645	82	874537	40.719	ng	100
26) 2-Nitrophenol	7.722	139	142080	40.062	ng	100
27) 2,4-Dimethylphenol	7.757	122	402429	41.112	ng	100
28) bis(2-Chloroethoxy)met...	7.857	93	535471	40.601	ng	100
29) 2,4-Dichlorophenol	7.957	162	372865	42.539	ng	100
30) 1,2,4-Trichlorobenzene	8.045	180	401673	40.473	ng	100
31) Naphthalene	8.128	128	1343175	40.888	ng	100
32) Benzoic acid	7.869	122	164746m	38.468	ng	
33) 4-Chloroaniline	8.175	127	595138	40.989	ng	100
34) Hexachlorobutadiene	8.245	225	225270	40.271	ng	100
35) Caprolactam	8.551	113	127357	43.662	ng	100
36) 4-Chloro-3-methylphenol	8.651	107	395374	42.445	ng	100
37) 2-Methylnaphthalene	8.822	142	889767	40.937	ng	100
38) 1-Methylnaphthalene	8.916	142	828740	40.806	ng	100
40) 1,2,4,5-Tetrachloroben...	8.980	216	391952	40.409	ng	100
41) Hexachlorocyclopentadiene	8.975	237	207179	43.318	ng	100
43) 2,4,6-Trichlorophenol	9.092	196	242723	42.755	ng	100

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102523\
 Data File : BF135940.D
 Acq On : 25 Oct 2023 13:35
 Operator : CG\JU
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICCC040

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 10/26/2023
 Supervised By :mohammad ahmed 10/26/2023

Quant Time: Oct 26 03:51:20 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 26 03:12:14 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.128	196	292071	43.331	ng	100
46) 1,1'-Biphenyl	9.286	154	1056553	40.602	ng	100
47) 2-Chloronaphthalene	9.310	162	790034	41.046	ng	100
48) 2-Nitroaniline	9.404	65	204768	40.650	ng	100
49) Acenaphthylene	9.727	152	1257300	41.406	ng	100
50) Dimethylphthalate	9.592	163	915321	41.100	ng	100
51) 2,6-Dinitrotoluene	9.645	165	182446	40.375	ng	100
52) Acenaphthene	9.898	154	798156	41.015	ng	100
53) 3-Nitroaniline	9.816	138	216678	40.054	ng	100
54) 2,4-Dinitrophenol	9.916	184	39309	40.241	ng	100
55) Dibenzofuran	10.069	168	1161468	41.314	ng	100
56) 4-Nitrophenol	9.969	139	165540	40.182	ng	100
57) 2,4-Dinitrotoluene	10.051	165	218120	39.360	ng	100
58) Fluorene	10.416	166	858822	40.833	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.186	232	223136	43.993	ng	100
60) Diethylphthalate	10.292	149	890692	41.940	ng	100
61) 4-Chlorophenyl-phenyle...	10.410	204	411904	40.599	ng	100
62) 4-Nitroaniline	10.433	138	213831	39.953	ng	100
63) Azobenzene	10.569	77	888073	41.206	ng	100
65) 4,6-Dinitro-2-methylph...	10.457	198	74239	40.010	ng	100
66) n-Nitrosodiphenylamine	10.527	169	791999	41.206	ng	100
67) 4-Bromophenyl-phenylether	10.898	248	272833	41.874	ng	100
68) Hexachlorobenzene	10.957	284	287313	41.279	ng	100
69) Atrazine	11.057	200	216830	41.383	ng	100
70) Pentachlorophenol	11.151	266	169202	40.542	ng	100
71) Phenanthrene	11.380	178	1336767	41.530	ng	100
72) Anthracene	11.433	178	1380601	41.889	ng	100
73) Carbazole	11.586	167	1232719	41.929	ng	100
74) Di-n-butylphthalate	11.927	149	1339240	42.807	ng	100
75) Fluoranthene	12.574	202	1411158	41.672	ng	100
77) Benzidine	12.698	184	391794	50.102	ng	100
78) Pyrene	12.804	202	1442670	42.367	ng	100
80) Butylbenzylphthalate	13.439	149	512423	40.471	ng	100
81) Benzo(a)anthracene	13.998	228	1150704	41.730	ng	100
82) 3,3'-Dichlorobenzidine	13.963	252	353816	42.174	ng	100
83) Chrysene	14.033	228	1124143	41.857	ng	100
84) Bis(2-ethylhexyl)phtha...	14.004	149	613791	43.835	ng	100
85) Di-n-octyl phthalate	14.621	149	920430	41.728	ng	100
87) Indeno(1,2,3-cd)pyrene	16.992	276	980377	44.534	ng	100
88) Benzo(b)fluoranthene	15.057	252	923172	39.956	ng	100
89) Benzo(k)fluoranthene	15.086	252	916031	40.368	ng	100
90) Benzo(a)pyrene	15.427	252	869653	40.840	ng	100
91) Dibenzo(a,h)anthracene	17.021	278	805494	42.777	ng	100
92) Benzo(g,h,i)perylene	17.456	276	813915	38.798	ng	100

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102523\
 Data File : BF135940.D
 Acq On : 25 Oct 2023 13:35
 Operator : CG\JU
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICCC040

Quant Time: Oct 26 03:51:20 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 26 03:12:14 2023
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Yogesh Patel 10/26/2023
 Supervised By :mohammad ahmed 10/26/2023

