

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062923\
 Data File : BF134028.D
 Acq On : 29 Jun 2023 10:18
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 06/30/2023
 Supervised By :mohammad ahmed 06/30/2023

Quant Time: Jun 29 12:47:45 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 26 23:12:04 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.851	152	114495	20.000	ng	0.00
21) Naphthalene-d8	8.128	136	400148	20.000	ng	0.00
39) Acenaphthene-d10	9.886	164	199802	20.000	ng	0.00
64) Phenanthrene-d10	11.380	188	383200	20.000	ng	0.00
76) Chrysene-d12	14.045	240	185030	20.000	ng	0.00
86) Perylene-d12	15.551	264	208680	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	537911	78.589	ng	0.00
7) Phenol-d6	6.492	99	646133	76.760	ng	0.00
23) Nitrobenzene-d5	7.416	82	618186	83.278	ng	0.00
42) 2,4,6-Tribromophenol	10.686	330	202668	80.902	ng	0.00
45) 2-Fluorobiphenyl	9.210	172	1116700	81.259	ng	0.00
79) Terphenyl-d14	12.980	244	997193	86.737	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.622	88	107506	38.092	ng	98
3) Pyridine	3.410	79	289491	39.380	ng	96
4) n-Nitrosodimethylamine	3.369	42	170591	40.631	ng	92
6) Aniline	6.516	93	403450	39.388	ng	97
8) 2-Chlorophenol	6.639	128	268573	38.654	ng	95
9) Benzaldehyde	6.404	77	183091	39.437	ng	98
10) Phenol	6.504	94	339457	38.355	ng	95
11) bis(2-Chloroethyl)ether	6.592	93	245983	37.751	ng	98
12) 1,3-Dichlorobenzene	6.792	146	291895	39.405	ng	96
13) 1,4-Dichlorobenzene	6.869	146	288620	38.575	ng	99
14) 1,2-Dichlorobenzene	7.022	146	274152	39.125	ng	98
15) Benzyl Alcohol	6.992	79	258151	39.782	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.128	45	424939	36.863	ng	99
17) 2-Methylphenol	7.104	107	237966	38.246	ng	98
18) Hexachloroethane	7.357	117	112692	40.621	ng	98
19) n-Nitroso-di-n-propyla...	7.263	70	200718	37.601	ng	97
20) 3+4-Methylphenols	7.257	107	290684	38.510	ng	91
22) Acetophenone	7.263	105	371504	40.823	ng	# 95
24) Nitrobenzene	7.434	77	312814	41.701	ng	94
25) Isophorone	7.669	82	535651	39.704	ng	99
26) 2-Nitrophenol	7.745	139	148182	41.179	ng	90
27) 2,4-Dimethylphenol	7.786	122	229635	40.314	ng	97
28) bis(2-Chloroethoxy)met...	7.881	93	311317	39.852	ng	98
29) 2,4-Dichlorophenol	7.986	162	228421	40.440	ng	95
30) 1,2,4-Trichlorobenzene	8.069	180	237161	40.692	ng	100
31) Naphthalene	8.151	128	778889	40.298	ng	99
32) Benzoic acid	7.898	122	166924m	39.108	ng	
33) 4-Chloroaniline	8.204	127	334092	40.408	ng	99
34) Hexachlorobutadiene	8.263	225	154845	42.118	ng	99
35) Caprolactam	8.580	113	70964	38.157	ng	96
36) 4-Chloro-3-methylphenol	8.686	107	256478	40.420	ng	99
37) 2-Methylnaphthalene	8.845	142	537386	39.316	ng	99
38) 1-Methylnaphthalene	8.945	142	505595	39.790	ng	98
40) 1,2,4,5-Tetrachloroben...	9.010	216	246160	41.585	ng	98
41) Hexachlorocyclopentadiene	8.992	237	101567	36.167	ng	99
43) 2,4,6-Trichlorophenol	9.122	196	163187	40.497	ng	100

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062923\
 Data File : BF134028.D
 Acq On : 29 Jun 2023 10:18
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Jun 29 12:47:45 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 26 23:12:04 2023
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 06/30/2023
 Supervised By :mohammad ahmed 06/30/2023

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.163	196	190163	40.119	ng	99
46) 1,1'-Biphenyl	9.310	154	642165	40.067	ng	98
47) 2-Chloronaphthalene	9.333	162	467026	39.826	ng	96
48) 2-Nitroaniline	9.433	65	176006	41.183	ng	98
49) Acenaphthylene	9.751	152	801779	40.025	ng	99
50) Dimethylphthalate	9.610	163	575695	39.693	ng	100
51) 2,6-Dinitrotoluene	9.675	165	126036	40.346	ng	# 85
52) Acenaphthene	9.922	154	466053	40.095	ng	98
53) 3-Nitroaniline	9.845	138	148405	39.600	ng	97
54) 2,4-Dinitrophenol	9.957	184	58322	34.631	ng	96
55) Dibenzofuran	10.098	168	717615	39.503	ng	99
56) 4-Nitrophenol	10.010	139	91410	34.871	ng	95
57) 2,4-Dinitrotoluene	10.086	165	163818	39.709	ng	100
58) Fluorene	10.439	166	561053	39.698	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.216	232	147694m	39.502	ng	
60) Diethylphthalate	10.310	149	591627	39.153	ng	99
61) 4-Chlorophenyl-phenyle...	10.433	204	266988	39.200	ng	99
62) 4-Nitroaniline	10.463	138	138519	36.676	ng	93
63) Azobenzene	10.592	77	563266	39.407	ng	96
65) 4,6-Dinitro-2-methylph...	10.498	198	84663	40.524	ng	97
66) n-Nitrosodiphenylamine	10.551	169	488195	39.874	ng	98
67) 4-Bromophenyl-phenylether	10.922	248	165880	40.727	ng	# 90
68) Hexachlorobenzene	10.992	284	175937	40.684	ng	96
69) Atrazine	11.080	200	126769	37.432	ng	97
70) Pentachlorophenol	11.186	266	101012	39.066	ng	98
71) Phenanthrene	11.410	178	760804	39.438	ng	99
72) Anthracene	11.463	178	786912	39.219	ng	99
73) Carbazole	11.621	167	688311	37.479	ng	99
74) Di-n-butylphthalate	11.945	149	851380	38.983	ng	99
75) Fluoranthene	12.604	202	755774	36.239	ng	99
77) Benzidine	12.727	184	178122	40.458	ng	99
78) Pyrene	12.833	202	736637	42.779	ng	99
80) Butylbenzylphthalate	13.457	149	274353	42.482	ng	100
81) Benzo(a)anthracene	14.033	228	504245	39.295	ng	99
82) 3,3'-Dichlorobenzidine	13.998	252	161687	39.771	ng	98
83) Chrysene	14.068	228	490047	39.428	ng	99
84) Bis(2-ethylhexyl)phtha...	14.021	149	344726	46.454	ng	99
85) Di-n-octyl phthalate	14.639	149	554138	50.820	ng	100
87) Indeno(1,2,3-cd)pyrene	17.103	276	605651	41.826	ng	98
88) Benzo(b)fluoranthene	15.104	252	443595	36.151	ng	99
89) Benzo(k)fluoranthene	15.133	252	482374	38.611	ng	98
90) Benzo(a)pyrene	15.486	252	469841	39.597	ng	99
91) Dibenzo(a,h)anthracene	17.121	278	499771	42.255	ng	99
92) Benzo(g,h,i)perylene	17.580	276	499280	40.935	ng	98

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062923\
 Data File : BF134028.D
 Acq On : 29 Jun 2023 10:18
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Jun 29 12:47:45 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 26 23:12:04 2023
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Yogesh Patel 06/30/2023
 Supervised By :mohammad ahmed 06/30/2023

