

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112423\
 Data File : BF136194.D
 Acq On : 24 Nov 2023 23:32
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 11/25/2023
 Supervised By :mohammad ahmed 11/27/2023

Quant Time: Nov 24 23:52:51 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 24 23:44:28 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.810	152	119266	20.000	ng	0.00
21) Naphthalene-d8	8.086	136	507155	20.000	ng	0.00
39) Acenaphthene-d10	9.845	164	262838	20.000	ng	0.00
64) Phenanthrene-d10	11.333	188	477948	20.000	ng	0.00
76) Chrysene-d12	13.980	240	224867	20.000	ng	0.00
86) Perylene-d12	15.462	264	260969	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	726390	96.975	ng	0.00
7) Phenol-d6	6.445	99	907913	98.502	ng	0.00
23) Nitrobenzene-d5	7.375	82	835296	91.909	ng	0.00
42) 2,4,6-Tribromophenol	10.633	330	258573	93.931	ng	0.00
45) 2-Fluorobiphenyl	9.163	172	1766687	99.358	ng	0.00
79) Terphenyl-d14	12.927	244	1678947	105.984	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.563	88	136971	45.664	ng	96
3) Pyridine	3.322	79	376798	46.752	ng	97
4) n-Nitrosodimethylamine	3.293	42	180911	48.938	ng	93
6) Aniline	6.475	93	573612	49.101	ng	96
8) 2-Chlorophenol	6.592	128	389239	48.272	ng	95
9) Benzaldehyde	6.363	77	256855	48.201	ng	96
10) Phenol	6.457	94	499187	49.591	ng	98
11) bis(2-Chloroethyl)ether	6.551	93	360068	48.670	ng	97
12) 1,3-Dichlorobenzene	6.751	146	412875	48.541	ng	96
13) 1,4-Dichlorobenzene	6.828	146	421786	49.454	ng	94
14) 1,2-Dichlorobenzene	6.981	146	404583	50.469	ng	94
15) Benzyl Alcohol	6.951	79	343330	49.804	ng	95
16) 2,2'-oxybis(1-Chloropr...	7.087	45	490566	50.757	ng	97
17) 2-Methylphenol	7.057	107	327571	48.689	ng	95
18) Hexachloroethane	7.316	117	145103	46.620	ng	94
19) n-Nitroso-di-n-propyla...	7.222	70	283465	53.276	ng	# 92
20) 3+4-Methylphenols	7.216	107	431259	51.364	ng	91
22) Acetophenone	7.216	105	579712	49.509	ng	# 88
24) Nitrobenzene	7.392	77	410090	46.366	ng	97
25) Isophorone	7.628	82	759129	48.401	ng	95
26) 2-Nitrophenol	7.704	139	173823	40.549	ng	98
27) 2,4-Dimethylphenol	7.739	122	316010	43.887	ng	92
28) bis(2-Chloroethoxy)met...	7.839	93	453097	47.903	ng	# 92
29) 2,4-Dichlorophenol	7.939	162	315617	44.954	ng	96
30) 1,2,4-Trichlorobenzene	8.028	180	358846	45.904	ng	94
31) Naphthalene	8.110	128	1206772	48.304	ng	96
32) Benzoic acid	7.857	122	201023m	38.848	ng	
33) 4-Chloroaniline	8.157	127	481610	46.478	ng	97
34) Hexachlorobutadiene	8.228	225	208448	43.864	ng	97
35) Caprolactam	8.528	113	91783	45.151	ng	# 78
36) 4-Chloro-3-methylphenol	8.639	107	349059	47.353	ng	99
37) 2-Methylnaphthalene	8.798	142	779038	46.508	ng	91
38) 1-Methylnaphthalene	8.898	142	724434	46.614	ng	89
40) 1,2,4,5-Tetrachloroben...	8.963	216	356573	45.266	ng	97
41) Hexachlorocyclopentadiene	8.951	237	177575	43.715	ng	98
43) 2,4,6-Trichlorophenol	9.075	196	221656	44.938	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.110	196	257900	45.387	ng	97
46) 1,1'-Biphenyl	9.263	154	987505	47.880	ng	96
47) 2-Chloronaphthalene	9.286	162	718442	47.657	ng	96
48) 2-Nitroaniline	9.386	65	209495	45.579	ng	88
49) Acenaphthylene	9.704	152	1136244	48.829	ng	96
50) Dimethylphthalate	9.569	163	857176	48.268	ng	98
51) 2,6-Dinitrotoluene	9.627	165	163606	42.456	ng	92
52) Acenaphthene	9.875	154	718094	46.332	ng	89
53) 3-Nitroaniline	9.798	138	183416	44.641	ng	95
54) 2,4-Dinitrophenol	9.904	184	63874	37.765	ng	# 85
55) Dibenzofuran	10.051	168	1047600	48.616	ng	98
56) 4-Nitrophenol	9.951	139	123236	45.137	ng	93
57) 2,4-Dinitrotoluene	10.033	165	218280	44.599	ng	93
58) Fluorene	10.392	166	798171	48.803	ng	92
59) 2,3,4,6-Tetrachlorophenol	10.169	232	193494	44.929	ng	94
60) Diethylphthalate	10.269	149	842737	47.303	ng	96
61) 4-Chlorophenyl-phenyle...	10.386	204	384876	46.293	ng	96
62) 4-Nitroaniline	10.416	138	171366	45.765	ng	96
63) Azobenzene	10.551	77	793525	50.906	ng	95
65) 4,6-Dinitro-2-methylph...	10.439	198	98229	36.913	ng	94
66) n-Nitrosodiphenylamine	10.510	169	691948	45.633	ng	93
67) 4-Bromophenyl-phenylether	10.880	248	237270	44.574	ng	92
68) Hexachlorobenzene	10.939	284	254649	45.060	ng	94
69) Atrazine	11.039	200	149866	39.013	ng	98
70) Pentachlorophenol	11.133	266	141229	44.682	ng	97
71) Phenanthrene	11.357	178	1170052	47.806	ng	96
72) Anthracene	11.410	178	1202692	48.140	ng	96
73) Carbazole	11.569	167	961076	48.008	ng	97
74) Di-n-butylphthalate	11.904	149	1220756	51.546	ng	# 96
75) Fluoranthene	12.551	202	1067741	47.772	ng	95
77) Benzidine	12.674	184	176381	45.299	ng	97
78) Pyrene	12.780	202	1040355	49.475	ng	95
80) Butylbenzylphthalate	13.410	149	333008	49.700	ng	97
81) Benzo(a)anthracene	13.974	228	731286	49.030	ng	96
82) 3,3'-Dichlorobenzidine	13.939	252	234641	49.004	ng	96
83) Chrysene	14.010	228	704550	48.284	ng	96
84) Bis(2-ethylhexyl)phtha...	13.974	149	429820	53.147	ng	96
85) Di-n-octyl phthalate	14.592	149	691048	51.036	ng	100
87) Indeno(1,2,3-cd)pyrene	16.968	276	838519	43.853	ng	99
88) Benzo(b)fluoranthene	15.027	252	711638	49.391	ng	95
89) Benzo(k)fluoranthene	15.057	252	661482	45.591	ng	97
90) Benzo(a)pyrene	15.398	252	671181	47.229	ng	98
91) Dibenzo(a,h)anthracene	16.992	278	680402	42.936	ng	99
92) Benzo(g,h,i)perylene	17.427	276	671775	41.812	ng	94

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

