

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082018\
 Data File : BF108280.D
 Acq On : 20 Aug 2018 10:51
 Operator : JU/SJ
 Sample : PB112114BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB112114BS

Manual Integrations
 APPROVED

Sohil
 8/21/2018 1:42:39 PM

Quant Time: Aug 20 11:34:34 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 08 12:24:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.02	152	163645	20.00	ng	0.00
21) Naphthalene-d8	8.30	136	677830	20.00	ng	0.00
38) Acenaphthene-d10	10.07	164	339621	20.00	ng	0.00
63) Phenanthrene-d10	11.56	188	680254	20.00	ng	0.00
75) Chrysene-d12	14.22	240	512778	20.00	ng	0.00
86) Perylene-d12	15.78	264	367667	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.66	112	1235054	136.64	ng	0.03
7) Phenol-d6	6.65	99	1640068	136.54	ng	0.02
23) Nitrobenzene-d5	7.58	82	1083360	89.19	ng	0.00
41) 2,4,6-Tribromophenol	10.87	330	495877	146.38	ng	0.00
44) 2-Fluorobiphenyl	9.38	172	2023392	95.65	ng	0.00
78) Terphenyl-d14	13.15	244	2282320	113.17	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.03	88	185899	40.026	ng	93
3) Pyridine	3.77	79	494979	41.372	ng	88
4) n-Nitrosodimethylamine	3.74	42	282106	41.904	ng	85
6) Aniline	6.68	93	630422	38.586	ng	97
8) 2-Chlorophenol	6.81	128	514079	50.498	ng	98
9) Benzaldehyde	6.57	77	228590	24.546	ng	98
10) Phenol	6.66	94	643810	51.818	ng	90
11) bis(2-Chloroethyl)ether	6.75	93	487956	48.579	ng	91
12) 1,3-Dichlorobenzene	6.96	146	574220	48.782	ng	99
13) 1,4-Dichlorobenzene	7.04	146	574540	48.580	ng	98
14) 1,2-Dichlorobenzene	7.19	146	533108	48.461	ng	97
15) Benzyl Alcohol	7.15	79	467049	46.189	ng	95
16) 2,2'-oxybis(1-Chloropropan	7.28	45	908163	43.888	ng	99
17) 2-Methylphenol	7.27	107	444293	50.892	ng	# 95
18) Hexachloroethane	7.53	117	210278	49.942	ng	91
19) n-Nitroso-di-n-propylamine	7.42	70	384454	47.332	ng	91
20) 3+4-Methylphenols	7.41	107	538040	48.093	ng	# 87
22) Acetophenone	7.42	105	727686	45.912	ng	# 93
24) Nitrobenzene	7.60	77	567044	46.163	ng	96
25) Isophorone	7.84	82	1055641	45.594	ng	96
26) 2-Nitrophenol	7.91	139	255536	55.087	ng	89
27) 2,4-Dimethylphenol	7.94	122	487214	47.413	ng	96
28) bis(2-Chloroethoxy)methane	8.04	93	635922	47.049	ng	99
29) 2,4-Dichlorophenol	8.15	162	460198	48.664	ng	99
30) 1,2,4-Trichlorobenzene	8.24	180	494268	47.406	ng	98
31) Naphthalene	8.32	128	1484160	48.146	ng	99
32) Benzoic acid	8.07	122	217290	37.878	ng	95
33) 4-Chloroaniline	8.37	127	415403	30.922	ng	98
34) Hexachlorobutadiene	8.43	225	314092	47.587	ng	99
35) Caprolactam	8.75	113	131741m	44.455	ng	
36) 4-Chloro-3-methylphenol	8.85	107	475007	46.562	ng	99
37) 2-Methylnaphthalene	9.02	142	1034703	47.442	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.18	216	513534	49.239	ng	98
40) Hexachlorocyclopentadiene	9.17	237	563991	83.722	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.30	196	338368	52.282	ng	99
43) 2,4,5-Trichlorophenol	9.34	196	366280	51.411	ng	96
45) 1,1'-Biphenyl	9.48	154	1234902	49.317	ng	98
46) 2-Chloronaphthalene	9.51	162	965359	48.778	ng	96
47) 2-Nitroaniline	9.61	65	319028	49.820	ng	93
48) Acenaphthylene	9.93	152	1532915	48.994	ng	99
49) Dimethylphthalate	9.78	163	1123045	47.471	ng	98
50) 2,6-Dinitrotoluene	9.85	165	250416	50.593	ng	96
51) Acenaphthene	10.10	154	900589	46.995	ng	99
52) 3-Nitroaniline	10.02	138	200289	36.985	ng	97
53) 2,4-Dinitrophenol	10.13	184	180507	95.916	ng #	20
54) Dibenzofuran	10.27	168	1330065	47.906	ng	96
55) 4-Nitrophenol	10.18	139	377067	91.948	ng	94
56) 2,4-Dinitrotoluene	10.25	165	335257	52.622	ng #	97
57) Fluorene	10.62	166	1066299	48.516	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.39	232	300948	50.539	ng #	85
59) Diethylphthalate	10.48	149	1080666	47.174	ng	96
60) 4-Chlorophenyl-phenylether	10.60	204	546392	47.710	ng	95
61) 4-Nitroaniline	10.64	138	265067	49.507	ng	95
62) Azobenzene	10.77	77	1103049	46.625	ng	94
64) 4,6-Dinitro-2-methylphenol	10.67	198	124257	48.581	ng	89
65) n-Nitrosodiphenylamine	10.72	169	958520	47.683	ng	97
66) 4-Bromophenyl-phenylether	11.10	248	359959	48.692	ng #	81
67) Hexachlorobenzene	11.17	284	372868	47.938	ng #	90
68) Atrazine	11.25	200	328707	48.752	ng	96
69) Pentachlorophenol	11.37	266	359125	96.463	ng	96
70) Phenanthrene	11.59	178	1540905	48.025	ng	99
71) Anthracene	11.64	178	1574920	47.892	ng	98
72) Carbazole	11.80	167	1390887	47.605	ng	99
73) Di-n-butylphthalate	12.11	149	1634343	49.385	ng	99
74) Fluoranthene	12.78	202	1656404	47.303	ng	96
76) Benzidine	12.90	184	793682	41.427	ng	98
77) Pyrene	13.02	202	1650231	50.832	ng	97
79) Butylbenzylphthalate	13.62	149	703655	54.585	ng #	97
80) Benzo(a)anthracene	14.21	228	1436452	48.812	ng	98
81) 3,3'-Dichlorobenzidine	14.17	252	351350	33.315	ng #	96
82) Chrysene	14.25	228	1308406	48.496	ng	98
83) Bis(2-ethylhexyl)phthalate	14.17	149	883030	50.584	ng	100
84) Di-n-octyl phthalate	14.79	149	1374879	51.642	ng	97
85) Indeno(1,2,3-cd)pyrene	17.41	276	992424	42.454	ng #	90
87) Benzo(b)fluoranthene	15.31	252	1145506	52.140	ng #	96
88) Benzo(k)fluoranthene	15.35	252	1006757	49.851	ng #	96
89) Benzo(a)pyrene	15.72	252	982612	49.947	ng	96
90) Dibenzo(a,h)anthracene	17.43	278	829406	50.954	ng #	94
91) Benzo(g,h,i)perylene	17.91	276	841796	52.519	ng #	90

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

