

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082818\
 Data File : BF108568.D
 Acq On : 28 Aug 2018 21:02
 Operator : JU/SJ
 Sample : J4535-04MSD
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RA-081618-FL-036MSD

Quant Time: Aug 29 01:36:29 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 22 11:01:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.00	152	104823	20.00	ng	0.00
21) Naphthalene-d8	8.28	136	397720	20.00	ng	0.00
38) Acenaphthene-d10	10.05	164	186860	20.00	ng	0.00
63) Phenanthrene-d10	11.54	188	314813	20.00	ng	0.00
75) Chrysene-d12	14.20	240	266673	20.00	ng	0.00
86) Perylene-d12	15.75	264	209930	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	546443	94.38	ng	0.01
7) Phenol-d6	6.64	99	719910	93.57	ng	0.00
23) Nitrobenzene-d5	7.57	82	473810	66.48	ng	0.00
41) 2,4,6-Tribromophenol	10.84	330	170770	91.62	ng	0.00
44) 2-Fluorobiphenyl	9.35	172	806843	69.32	ng	-0.01
78) Terphenyl-d14	13.12	244	719839	68.63	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.90	88	67070	22.544	ng	93
3) Pyridine	3.70	79	178455	23.286	ng	87
4) n-Nitrosodimethylamine	3.65	42	106552	24.709	ng	# 79
6) Aniline	6.67	93	253884	24.259	ng	# 85
8) 2-Chlorophenol	6.79	128	210873	32.338	ng	91
9) Benzaldehyde	6.56	77	102482	17.180	ng	98
10) Phenol	6.65	94	262176	32.943	ng	82
11) bis(2-Chloroethyl)ether	6.73	93	201995	31.394	ng	91
12) 1,3-Dichlorobenzene	6.94	146	234098	31.048	ng	99
13) 1,4-Dichlorobenzene	7.02	146	240873	31.796	ng	98
14) 1,2-Dichlorobenzene	7.17	146	225351	31.981	ng	97
15) Benzyl Alcohol	7.14	79	191150	29.512	ng	93
16) 2,2'-oxybis(1-Chloropropan	7.27	45	367160	27.700	ng	82
17) 2-Methylphenol	7.25	107	173487	31.024	ng	# 89
18) Hexachloroethane	7.51	117	77872	28.873	ng	94
19) n-Nitroso-di-n-propylamine	7.40	70	157252	30.224	ng	89
20) 3+4-Methylphenols	7.41	107	225977	31.534	ng	# 40
22) Acetophenone	7.41	105	307066	33.019	ng	# 88
24) Nitrobenzene	7.58	77	232037	32.195	ng	95
25) Isophorone	7.81	82	414961	30.545	ng	96
26) 2-Nitrophenol	7.90	139	97984	35.999	ng	95
27) 2,4-Dimethylphenol	7.93	122	175557	29.116	ng	97
28) bis(2-Chloroethoxy)methane	8.02	93	250999	31.649	ng	98
29) 2,4-Dichlorophenol	8.14	162	184367	33.227	ng	99
30) 1,2,4-Trichlorobenzene	8.22	180	194397	31.776	ng	95
31) Naphthalene	8.31	128	608774	33.657	ng	100
32) Benzoic acid	8.01	122	21288	11.301	ng	94
33) 4-Chloroaniline	8.35	127	203538	25.822	ng	96
34) Hexachlorobutadiene	8.41	225	126870	32.759	ng	99
35) Caprolactam	8.71	113	51312	29.509	ng	# 75
36) 4-Chloro-3-methylphenol	8.84	107	187589	31.339	ng	96
37) 2-Methylnaphthalene	9.00	142	411970	32.192	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.16	216	203892	35.532	ng	# 96
40) Hexachlorocyclopentadiene	9.14	237	38163	10.296	ng	97

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42) 2,4,6-Trichlorophenol	9.28	196	129696	36.422	nq	100
43) 2,4,5-Trichlorophenol	9.32	196	134915	34.418	nq	93
45) 1,1'-Biphenyl	9.46	154	494215	35.872	nq	98
46) 2-Chloronaphthalene	9.49	162	375405	34.476	nq	99
47) 2-Nitroaniline	9.59	65	123932	35.175	nq	85
48) Acenaphthylene	9.91	152	583439	33.892	nq	100
49) Dimethylphthalate	9.75	163	502600	38.613	nq	99
50) 2,6-Dinitrotoluene	9.82	165	92293	33.890	nq	90
51) Acenaphthene	10.08	154	330979	31.391	nq	98
52) 3-Nitroaniline	10.01	138	87680	29.427	nq	100
53) 2,4-Dinitrophenol	10.11	184	10762	16.654	nq #	23
54) Dibenzofuran	10.25	168	498307	32.621	nq	98
55) 4-Nitrophenol	10.17	139	109323	48.452	nq #	79
56) 2,4-Dinitrotoluene	10.24	165	115987	33.088	nq	98
57) Fluorene	10.60	166	392845	32.486	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.37	232	96859	29.563	nq #	90
59) Diethylphthalate	10.45	149	426980	33.876	nq	97
60) 4-Chlorophenyl-phenylether	10.58	204	199476	31.657	nq	92
61) 4-Nitroaniline	10.62	138	83348	28.293	nq	95
62) Azobenzene	10.74	77	396257	30.442	nq	94
64) 4,6-Dinitro-2-methylphenol	10.64	198	13196	15.277	nq	91
65) n-Nitrosodiphenylamine	10.70	169	347406	37.343	nq	96
66) 4-Bromophenyl-phenylether	11.08	248	121045	35.381	nq #	84
67) Hexachlorobenzene	11.15	284	121243	33.682	nq #	84
68) Atrazine	11.22	200	112154	35.944	nq	95
69) Pentachlorophenol	11.35	266	98658	57.262	nq	100
70) Phenanthrene	11.57	178	504075	33.948	nq	99
71) Anthracene	11.62	178	513485	33.740	nq	100
72) Carbazole	11.78	167	442828	32.750	nq	99
73) Di-n-butylphthalate	12.08	149	559829	36.553	nq	100
74) Fluoranthene	12.76	202	520319	32.108	nq	95
76) Benzidine	12.88	184	272883	27.388	nq	98
77) Pyrene	12.99	202	515751	30.548	nq	100
79) Butylbenzylphthalate	13.59	149	218424	32.581	nq	96
80) Benzo(a)anthracene	14.19	228	497841	32.529	nq	99
81) 3,3'-Dichlorobenzidine	14.14	252	181861	33.158	nq #	98
82) Chrysene	14.22	228	463706	33.049	nq	100
83) Bis(2-ethylhexyl)phthalate	14.14	149	301051	33.161	nq	99
84) Di-n-octyl phthalate	14.77	149	530236	38.296	nq	96
85) Indeno(1,2,3-cd)pyrene	17.37	276	334083	27.480	nq #	89
87) Benzo(b)fluoranthene	15.28	252	425026	33.882	nq #	96
88) Benzo(k)fluoranthene	15.32	252	392683	34.054	nq #	94
89) Benzo(a)pyrene	15.69	252	367472	32.714	nq #	96
90) Dibenzo(a,h)anthracene	17.38	278	288016	30.989	nq #	92
91) Benzo(g,h,i)perylene	17.87	276	264539	28.905	nq #	89

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

