

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
 Data File : BF102504.D
 Acq On : 27 Jan 2018 12:23
 Operator : SJ/JU
 Sample : J1306-01MSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC1MSD

Manual Integrations
 APPROVED

Sohil
 1/29/2018 5:44:24 PM

Quant Time: Jan 27 21:18:26 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jan 27 20:48:40 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.71	152	180587	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	761898	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	344897	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	584717	20.00	ng	0.00
75) Chrysene-d12	13.87	240	318570	20.00	ng	0.00
86) Perylene-d12	15.30	264	294139	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.33	112	893680	75.04	ng	0.01
7) Phenol-d6	6.36	99	1086612	75.68	ng	0.00
23) Nitrobenzene-d5	7.28	82	679698	51.27	ng	0.00
41) 2,4,6-Tribromophenol	10.55	330	202632	59.29	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	1033533	44.06	ng	0.00
78) Terphenyl-d14	12.82	244	843934	59.72	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.43	88	134319	23.736	ng	99
3) Pyridine	3.14	79	352833	23.859	ng	96
4) n-Nitrosodimethylamine	3.10	42	170192	29.356	ng	97
6) Aniline	6.38	93	334524	17.614	ng	# 64
8) 2-Chlorophenol	6.50	128	361108	28.924	ng	95
9) Benzaldehyde	6.26	77	171876	18.208	ng	93
10) Phenol	6.37	94	451086	28.776	ng	77
11) bis(2-Chloroethyl)ether	6.45	93	354612	29.743	ng	97
12) 1,3-Dichlorobenzene	6.65	146	373605	25.162	ng	99
13) 1,4-Dichlorobenzene	6.73	146	384607	25.858	ng	100
14) 1,2-Dichlorobenzene	6.88	146	358308	26.337	ng	99
15) Benzyl Alcohol	6.86	79	286858	28.315	ng	99
16) 2,2'-oxybis(1-Chloropropan	6.99	45	547148	27.363	ng	91
17) 2-Methylphenol	6.98	107	295623	27.264	ng	# 92
18) Hexachloroethane	7.22	117	134390	25.929	ng	91
19) n-Nitroso-di-n-propylamine	7.13	70	245978	27.993	ng	96
20) 3+4-Methylphenols	7.13	107	367725	28.836	ng	# 86
22) Acetophenone	7.13	105	492715	27.128	ng	# 93
24) Nitrobenzene	7.30	77	368177	27.136	ng	100
25) Isophorone	7.54	82	683198	28.905	ng	99
26) 2-Nitrophenol	7.62	139	206316	28.892	ng	96
27) 2,4-Dimethylphenol	7.66	122	321413	30.997	ng	97
28) bis(2-Chloroethoxy)methane	7.75	93	437842	29.988	ng	99
29) 2,4-Dichlorophenol	7.86	162	303005	28.688	ng	96
30) 1,2,4-Trichlorobenzene	7.94	180	283109	25.005	ng	99
31) Naphthalene	8.02	128	988400	25.983	ng	99
32) Benzoic acid	7.75	122	29204	3.362	ng	95
33) 4-Chloroaniline	8.07	127	259647	16.232	ng	97
34) Hexachlorobutadiene	8.13	225	148436	24.547	ng	98
35) Caprolactam	8.44	113	91649m	26.273	ng	
36) 4-Chloro-3-methylphenol	8.56	107	322963	29.009	ng	100
37) 2-Methylnaphthalene	8.71	142	672793	26.557	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.87	216	264604	25.500	ng	99
40) Hexachlorocyclopentadiene	8.86	237	135175	29.109	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.99	196	188896	27.194	ng	99
43) 2,4,5-Trichlorophenol	9.04	196	195662	25.017	ng	96
45) 1,1'-Biphenyl	9.17	154	779570	25.278	ng	98
46) 2-Chloronaphthalene	9.20	162	601861	25.108	ng	98
47) 2-Nitroaniline	9.30	65	206300	27.605	ng	98
48) Acenaphthylene	9.62	152	915732	24.521	ng	99
49) Dimethylphthalate	9.47	163	854333	29.968	ng	100
50) 2,6-Dinitrotoluene	9.55	165	162655	26.464	ng	95
51) Acenaphthene	9.79	154	536888	24.616	ng	99
52) 3-Nitroaniline	9.72	138	147817	20.330	ng	93
53) 2,4-Dinitrophenol	9.83	184	51284	21.433	ng	# 20
54) Dibenzofuran	9.96	168	780710	26.448	ng	98
55) 4-Nitrophenol	9.89	139	213641	44.512	ng	94
56) 2,4-Dinitrotoluene	9.95	165	198106	26.210	ng	# 92
57) Fluorene	10.30	166	624607	26.704	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.08	232	152031	27.225	ng	98
59) Diethylphthalate	10.17	149	703807	25.406	ng	98
60) 4-Chlorophenyl-phenylether	10.29	204	274487	27.067	ng	98
61) 4-Nitroaniline	10.33	138	175332	24.166	ng	88
62) Azobenzene	10.45	77	658435	25.960	ng	97
64) 4,6-Dinitro-2-methylphenol	10.36	198	74629	19.131	ng	84
65) n-Nitrosodiphenylamine	10.41	169	575296	26.806	ng	100
66) 4-Bromophenyl-phenylether	10.78	248	167585	26.624	ng	92
67) Hexachlorobenzene	10.85	284	164844	23.417	ng	97
68) Atrazine	10.95	200	181219	28.590	ng	99
69) Pentachlorophenol	11.05	266	138343	37.410	ng	100
70) Phenanthrene	11.26	178	863975	25.357	ng	99
71) Anthracene	11.32	178	894058	25.708	ng	99
72) Carbazole	11.47	167	831675	24.513	ng	99
73) Di-n-butylphthalate	11.80	149	1105817	26.075	ng	99
74) Fluoranthene	12.45	202	829630	23.877	ng	97
76) Benzidine	12.57	184	186989	17.472	ng	96
77) Pyrene	12.68	202	814037	33.050	ng	99
79) Butylbenzylphthalate	13.29	149	404746	33.020	ng	97
80) Benzo(a)anthracene	13.86	228	545329	27.804	ng	99
81) 3,3'-Dichlorobenzidine	13.83	252	139002	19.320	ng	98
82) Chrysene	13.90	228	504520	25.331	ng	99
83) Bis(2-ethylhexyl)phthalate	13.84	149	560075	34.000	ng	# 98
84) Di-n-octyl phthalate	14.46	149	780802	29.146	ng	100
85) Indeno(1,2,3-cd)pyrene	16.70	276	515334	31.329	ng	99
87) Benzo(b)fluoranthene	14.89	252	441650	23.166	ng	98
88) Benzo(k)fluoranthene	14.92	252	441285	24.500	ng	99
89) Benzo(a)pyrene	15.24	252	438984	25.531	ng	98
90) Dibenzo(a,h)anthracene	16.71	278	427471	30.691	ng	99
91) Benzo(g,h,i)perylene	17.12	276	441250	32.085	ng	99

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