

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF022818\
 Data File : BF103311.D
 Acq On : 28 Feb 2018 23:24
 Operator : SJ/JU
 Sample : J1708-01MSD
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RO-222-226MSD

Manual Integrations
 APPROVED

Sohil
 3/1/2018 11:43:31 AM

Quant Time: Mar 01 01:25:15 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 27 17:05:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	144988	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	616752	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	265262	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	411214	20.00	ng	0.00
75) Chrysene-d12	14.06	240	240941	20.00	ng	0.00
86) Perylene-d12	15.56	264	241568	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.50	112	1229212	128.22	ng	0.01
7) Phenol-d6	6.52	99	1496932	127.61	ng	0.00
23) Nitrobenzene-d5	7.45	82	912244	84.47	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	245646	109.40	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1298607	84.11	ng	0.00
78) Terphenyl-d14	12.99	244	950668	94.85	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.70	88	204192	40.329	ng	93
3) Pyridine	3.49	79	530411	39.240	ng	97
4) n-Nitrosodimethylamine	3.43	42	320939	58.872	ng	# 90
6) Aniline	6.55	93	323506	19.109	ng	94
8) 2-Chlorophenol	6.66	128	438051	43.987	ng	91
9) Benzaldehyde	6.43	77	228149	29.708	ng	95
10) Phenol	6.53	94	597831	43.901	ng	88
11) bis(2-Chloroethyl)ether	6.61	93	469870	45.462	ng	91
12) 1,3-Dichlorobenzene	6.82	146	449074	39.460	ng	98
13) 1,4-Dichlorobenzene	6.89	146	448481	39.306	ng	98
14) 1,2-Dichlorobenzene	7.05	146	401918	38.202	ng	99
15) Benzyl Alcohol	7.02	79	373949	42.304	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.15	45	703188	39.620	ng	91
17) 2-Methylphenol	7.13	107	370713	40.962	ng	94
18) Hexachloroethane	7.38	117	149818	36.069	ng	91
19) n-Nitroso-di-n-propylamine	7.29	70	295893	40.530	ng	97
20) 3+4-Methylphenols	7.29	107	436003	39.522	ng	# 57
22) Acetophenone	7.29	105	560466	38.932	ng	# 89
24) Nitrobenzene	7.46	77	479606	40.336	ng	99
25) Isophorone	7.70	82	893684	42.016	ng	98
26) 2-Nitrophenol	7.77	139	229465	40.994	ng	96
27) 2,4-Dimethylphenol	7.81	122	379089	44.306	ng	99
28) bis(2-Chloroethoxy)methane	7.90	93	559904	44.772	ng	99
29) 2,4-Dichlorophenol	8.02	162	353836	43.195	ng	98
30) 1,2,4-Trichlorobenzene	8.10	180	331432	37.994	ng	98
31) Naphthalene	8.18	128	1186765	39.303	ng	100
32) Benzoic acid	7.93	122	92169	19.342	ng	99
33) 4-Chloroaniline	8.23	127	121931	9.358	ng	95
34) Hexachlorobutadiene	8.29	225	164989	36.321	ng	97
35) Caprolactam	8.61	113	124467m	43.233	ng	
36) 4-Chloro-3-methylphenol	8.72	107	391475	42.369	ng	98
37) 2-Methylnaphthalene	8.87	142	773801	39.653	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.04	216	285985	39.437	ng	99
40) Hexachlorocyclopentadiene	9.02	237	47689	25.467	ng	99

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42) 2,4,6-Trichlorophenol	9.15	196	208858	42.803	nq	98
43) 2,4,5-Trichlorophenol	9.20	196	221474	39.463	nq	96
45) 1,1'-Biphenyl	9.33	154	864964	38.914	nq	98
46) 2-Chloronaphthalene	9.36	162	692194	39.363	nq	99
47) 2-Nitroaniline	9.46	65	285708	43.862	nq	89
48) Acenaphthylene	9.78	152	1043690	38.575	nq	100
49) Dimethylphthalate	9.63	163	953061	44.787	nq	100
50) 2,6-Dinitrotoluene	9.70	165	181272	40.593	nq #	87
51) Acenaphthene	9.95	154	602727	38.203	nq	99
52) 3-Nitroaniline	9.87	138	133276	23.524	nq	99
53) 2,4-Dinitrophenol	9.99	184	28019	24.902	nq #	19
54) Dibenzofuran	10.12	168	873383	40.327	nq	97
55) 4-Nitrophenol	10.05	139	238736	67.989	nq	89
56) 2,4-Dinitrotoluene	10.12	165	221891	39.288	nq #	92
57) Fluorene	10.47	166	685899	37.177	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.25	232	153140	40.620	nq	98
59) Diethylphthalate	10.33	149	808106	39.706	nq	99
60) 4-Chlorophenyl-phenylether	10.45	204	313167	40.028	nq	99
61) 4-Nitroaniline	10.49	138	191498	33.840	nq	99
62) Azobenzene	10.62	77	827965	39.967	nq	98
64) 4,6-Dinitro-2-methylphenol	10.53	198	35626	17.717	nq	94
65) n-Nitrosodiphenylamine	10.57	169	627950	43.858	nq	100
66) 4-Bromophenyl-phenylether	10.95	248	183614	42.864	nq	93
67) Hexachlorobenzene	11.02	284	180849	38.897	nq	97
68) Atrazine	11.10	200	188627	43.831	nq	100
69) Pentachlorophenol	11.22	266	138739	61.714	nq	97
70) Phenanthrene	11.43	178	862570	38.116	nq	99
71) Anthracene	11.49	178	913273	39.355	nq	99
72) Carbazole	11.64	167	874753	37.853	nq	99
73) Di-n-butylphthalate	11.96	149	1185126	40.984	nq	99
74) Fluoranthene	12.63	202	782802	34.422	nq	98
76) Benzidine	12.74	184	399035	44.300	nq	98
77) Pyrene	12.86	202	772853	41.142	nq	99
79) Butylbenzylphthalate	13.46	149	442376	44.572	nq	95
80) Benzo(a)anthracene	14.04	228	628632	40.211	nq	99
81) 3,3'-Dichlorobenzidine	14.00	252	215397	38.608	nq	99
82) Chrysene	14.09	228	599001	39.461	nq	99
83) Bis(2-ethylhexyl)phthalate	14.02	149	577824	43.142	nq	98
84) Di-n-octyl phthalate	14.64	149	961083	44.413	nq	97
85) Indeno(1,2,3-cd)pyrene	17.09	276	539235	44.872	nq	97
87) Benzo(b)fluoranthene	15.12	252	616849	38.837	nq	98
88) Benzo(k)fluoranthene	15.15	252	582649	38.957	nq	99
89) Benzo(a)pyrene	15.50	252	592335	41.762	nq	98
90) Dibenzo(a,h)anthracene	17.10	278	456142	41.620	nq	99
91) Benzo(g,h,i)perylene	17.55	276	439029	40.987	nq	98

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

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