

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120718\
 Data File : BF111200.D
 Acq On : 7 Dec 2018 21:22
 Operator : JU/SJ
 Sample : J6214-18MSD
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-10(13-15)MSD

Manual Integrations
 APPROVED

Sohil
 12/11/2018 8:44:48 AM

Quant Time: Dec 08 02:04:04 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 05 17:30:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	478265	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	1917663	20.00	ng	0.00
39) Acenaphthene-d10	9.85	164	696094	20.00	ng	0.00
64) Phenanthrene-d10	11.33	188	1242692	20.00	ng	0.00
76) Chrysene-d12	13.96	240	760361	20.00	ng	0.01
87) Perylene-d12	15.40	264	801989	20.00	ng	0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.43	112	3720740	121.40	ng	0.01
7) Phenol-d6	6.45	99	4834896	122.46	ng	0.01
23) Nitrobenzene-d5	7.37	82	3121059	96.36	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	795781	143.82	ng	0.00
45) 2-Fluorobiphenyl	9.17	172	3451108	93.15	ng	0.00
79) Terphenyl-d14	12.91	244	2899103	78.09	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.56	88	566050	32.108	ng	96
3) Pyridine	3.30	79	1625360	41.570	ng	90
4) n-Nitrosodimethylamine	3.23	42	828294	43.026	ng	84
6) Aniline	6.47	93	844691	16.014	ng	97
8) 2-Chlorophenol	6.59	128	1442002	40.819	ng	94
9) Benzaldehyde	6.35	77	742691	33.432	ng	99
10) Phenol	6.46	94	2037666	47.334	ng	98
11) bis(2-Chloroethyl)ether	6.55	93	1429852	40.639	ng	96
12) 1,3-Dichlorobenzene	6.75	146	1480571	39.408	ng	95
13) 1,4-Dichlorobenzene	6.82	146	1490111	39.412	ng	98
14) 1,2-Dichlorobenzene	6.98	146	1397186	39.726	ng	100
15) Benzyl Alcohol	6.95	79	1303450	41.667	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.09	45	2704426	39.514	ng	98
17) 2-Methylphenol	7.06	107	1250321	43.116	ng	98
18) Hexachloroethane	7.32	117	565604	39.878	ng	96
19) n-Nitroso-di-n-propylamine	7.22	70	1060703	39.978	ng	93
20) 3+4-Methylphenols	7.21	107	1449889	40.994	ng	96
22) Acetophenone	7.22	105	1896975	40.934	ng	95
24) Nitrobenzene	7.39	77	1554714	44.615	ng	98
25) Isophorone	7.63	82	2833372	47.664	ng	100
26) 2-Nitrophenol	7.70	139	778808	52.894	ng	90
27) 2,4-Dimethylphenol	7.75	122	1330833	47.993	ng	98
28) bis(2-Chloroethoxy)methane	7.84	93	1832365	45.533	ng	98
29) 2,4-Dichlorophenol	7.95	162	1204746	44.131	ng	98
30) 1,2,4-Trichlorobenzene	8.03	180	1155897	41.949	ng	98
31) Naphthalene	8.11	128	3869424	46.344	ng	99
32) Benzoic acid	7.80	122	159588	8.727	ng	98
33) 4-Chloroaniline	8.15	127	230162	5.703	ng	96
34) Hexachlorobutadiene	8.23	225	602295	40.222	ng	98
35) Caprolactam	8.52	113	440492m	43.947	ng	
36) 4-Chloro-3-methylphenol	8.64	107	1061003	35.482	ng	99
37) 2-Methylnaphthalene	8.80	142	2061616	35.797	ng	98
38) 1-Methylnaphthalene	8.90	142	1959239	35.384	ng	97
40) 1,2,4,5-Tetrachlorobenzene	8.97	216	749627	40.406	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.96	237	813975	83.208	nq	98
43) 2,4,6-Trichlorophenol	9.08	196	606554	44.765	nq	100
44) 2,4,5-Trichlorophenol	9.12	196	641414	46.697	nq	99
46) 1,1'-Biphenyl	9.27	154	2329893	40.887	nq	97
47) 2-Chloronaphthalene	9.29	162	1876154	42.370	nq	96
48) 2-Nitroaniline	9.38	65	660932	46.058	nq	94
49) Acenaphthylene	9.70	152	2947848	47.627	nq	99
50) Dimethylphthalate	9.57	163	2638501	54.463	nq	99
51) 2,6-Dinitrotoluene	9.62	165	511835	49.832	nq	90
52) Acenaphthene	9.88	154	1871585	45.771	nq	99
53) 3-Nitroaniline	9.79	138	228354	17.699	nq	91
54) 2,4-Dinitrophenol	9.89	184	309380	93.821	nq	# 74
55) Dibenzofuran	10.05	168	2591662	46.016	nq	95
56) 4-Nitrophenol	9.95	139	857024	88.151	nq	96
57) 2,4-Dinitrotoluene	10.03	165	680152	51.273	nq	97
58) Fluorene	10.39	166	1929853	48.591	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.17	232	514465m	50.184	nq	
60) Diethylphthalate	10.27	149	2160152	44.618	nq	98
61) 4-Chlorophenyl-phenylether	10.39	204	870564	44.479	nq	92
62) 4-Nitroaniline	10.40	138	497427	42.718	nq	95
63) Azobenzene	10.55	77	2195156	42.588	nq	96
65) 4,6-Dinitro-2-methylphenol	10.43	198	263866	51.396	nq	91
66) n-Nitrosodiphenylamine	10.50	169	1794848	41.248	nq	99
67) 4-Bromophenyl-phenylether	10.87	248	549470	41.770	nq	90
68) Hexachlorobenzene	10.94	284	576932	42.892	nq	95
69) Atrazine	11.03	200	579730	Below Cal		96
70) Pentachlorophenol	11.13	266	669500	75.688	nq	99
71) Phenanthrene	11.35	178	2816031	47.110	nq	100
72) Anthracene	11.40	178	2827816	46.932	nq	98
73) Carbazole	11.55	167	2619457	41.780	nq	98
74) Di-n-butylphthalate	11.89	149	3242691	44.184	nq	99
75) Fluoranthene	12.53	202	2687368	45.944	nq	98
77) Benzidine	12.65	184	559353	19.338	nq	95
78) Pyrene	12.76	202	2619078	44.213	nq	99
80) Butylbenzylphthalate	13.39	149	1501872	52.482	nq	99
81) Benzo(a)anthracene	13.95	228	1964921	44.106	nq	99
82) 3,3'-Dichlorobenzidine	13.91	252	229237	13.635	nq	97
83) Chrysene	13.99	228	2039315	46.111	nq	98
84) Bis(2-ethylhexyl)phthalate	13.95	149	1673546	48.416	nq	99
85) Di-n-octyl phthalate	14.56	149	2484894	45.729	nq	97
86) Indeno(1,2,3-cd)pyrene	16.82	276	2189364	61.656	nq	# 88
88) Benzo(b)fluoranthene	14.99	252	1991050	43.813	nq	98
89) Benzo(k)fluoranthene	15.02	252	1818770	41.399	nq	99
90) Benzo(a)pyrene	15.34	252	1918673	46.030	nq	97
91) Dibenzo(a,h)anthracene	16.83	278	1807340	51.668	nq	99
92) Benzo(g,h,i)perylene	17.24	276	2240296	63.733	nq	98

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

