

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040518\
 Data File : BF104278.D
 Acq On : 5 Apr 2018 19:15
 Operator : JU/SJ
 Sample : J2183-06MS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ES-4-DWS-1-NW@4.5MS

Manual Integrations
 APPROVED

Sohil
 4/6/2018 11:57:52 AM

Quant Time: Apr 06 01:33:04 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 04 18:06:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	128574	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	562306	20.00	ng	0.00
38) Acenaphthene-d10	9.99	164	260511	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	478462	20.00	ng	0.00
75) Chrysene-d12	14.14	240	356917	20.00	ng	0.00
86) Perylene-d12	15.67	264	309582	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.59	112	1035671	128.46	ng	0.00
7) Phenol-d6	6.59	99	1288526	129.90	ng	-0.01
23) Nitrobenzene-d5	7.52	82	826161	89.85	ng	0.00
41) 2,4,6-Tribromophenol	10.79	330	372277	128.34	ng	0.00
44) 2-Fluorobiphenyl	9.30	172	1416913	85.77	ng	0.00
78) Terphenyl-d14	13.06	244	1406741	91.00	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.85	88	126419	36.356	ng	# 85
3) Pyridine	3.67	79	241624	27.448	ng	# 82
4) n-Nitrosodimethylamine	3.64	42	55257	21.118	ng	# 77
6) Aniline	6.63	93	270080	22.745	ng	# 75
8) 2-Chlorophenol	6.74	128	443573	50.156	ng	97
9) Benzaldehyde	6.53	77	53361	7.902	ng	97
10) Phenol	6.60	94	626642	57.017	ng	95
11) bis(2-Chloroethyl)ether	6.69	93	455132	48.057	ng	90
12) 1,3-Dichlorobenzene	6.89	146	417847	42.031	ng	98
13) 1,4-Dichlorobenzene	6.96	146	497951	46.746	ng	98
14) 1,2-Dichlorobenzene	7.12	146	429364	44.845	ng	98
15) Benzyl Alcohol	7.10	79	270348m	41.919	ng	
16) 2,2'-oxybis(1-Chloropropan	7.21	45	474859	46.023	ng	# 45
17) 2-Methylphenol	7.20	107	338073	46.967	ng	# 78
18) Hexachloroethane	7.45	117	158763	44.515	ng	99
19) n-Nitroso-di-n-propylamine	7.36	70	288295	48.975	ng	89
20) 3+4-Methylphenols	7.36	107	405489	44.139	ng	# 51
22) Acetophenone	7.36	105	596577	43.850	ng	99
24) Nitrobenzene	7.54	77	378786	39.154	ng	96
25) Isophorone	7.76	82	829711	48.965	ng	99
26) 2-Nitrophenol	7.85	139	236570	46.846	ng	93
27) 2,4-Dimethylphenol	7.88	122	371043	50.946	ng	98
28) bis(2-Chloroethoxy)methane	7.97	93	526621	49.587	ng	100
29) 2,4-Dichlorophenol	8.09	162	363811	47.825	ng	98
30) 1,2,4-Trichlorobenzene	8.17	180	376206	43.587	ng	99
31) Naphthalene	8.25	128	1313914	47.830	ng	100
32) Benzoic acid	8.02	122	82745	15.509	ng	# 84
33) 4-Chloroaniline	8.33	127	123864	10.695	ng	93
34) Hexachlorobutadiene	8.35	225	190447	40.548	ng	100
35) Caprolactam	8.69	113	93532m	38.777	ng	
36) 4-Chloro-3-methylphenol	8.80	107	385212	47.875	ng	97
37) 2-Methylnaphthalene	8.94	142	808992	44.708	ng	100
39) 1,2,4,5-Tetrachlorobenzene	9.10	216	368723	46.156	ng	98
40) Hexachlorocyclopentadiene	9.08	237	239504	64.791	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.22	196	217639	43.024	nq	99
43) 2,4,5-Trichlorophenol	9.27	196	272124	44.553	nq	97
45) 1,1'-Biphenyl	9.40	154	1004378	44.921	nq	99
46) 2-Chloronaphthalene	9.43	162	800834	45.407	nq	99
47) 2-Nitroaniline	9.55	65	225606	44.860	nq	98
48) Acenaphthylene	9.85	152	1136950	42.551	nq	99
49) Dimethylphthalate	9.70	163	1050263	51.063	nq	99
50) 2,6-Dinitrotoluene	9.78	165	201877	45.621	nq	93
51) Acenaphthene	10.02	154	645812	41.781	nq	99
52) 3-Nitroaniline	9.97	138	120057	23.602	nq	# 96
53) 2,4-Dinitrophenol	10.07	184	96672	52.128	nq	# 38
54) Dibenzofuran	10.20	168	980863	43.455	nq	96
55) 4-Nitrophenol	10.15	139	181127	59.385	nq	95
56) 2,4-Dinitrotoluene	10.19	165	256790	45.477	nq	# 89
57) Fluorene	10.54	166	813738	43.704	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.32	232	247233	54.028	nq	90
59) Diethylphthalate	10.40	149	855163	43.400	nq	98
60) 4-Chlorophenyl-phenylether	10.52	204	382983	44.785	nq	99
61) 4-Nitroaniline	10.60	138	183866	38.630	nq	94
62) Azobenzene	10.69	77	796800	44.721	nq	94
64) 4,6-Dinitro-2-methylphenol	10.60	198	117689	40.635	nq	81
65) n-Nitrosodiphenylamine	10.65	169	723687	44.661	nq	99
66) 4-Bromophenyl-phenylether	11.02	248	239506	46.789	nq	93
67) Hexachlorobenzene	11.09	284	258488	43.899	nq	# 89
68) Atrazine	11.17	200	228024	45.934	nq	98
69) Pentachlorophenol	11.29	266	242474	74.834	nq	96
70) Phenanthrene	11.51	178	1178802	45.506	nq	99
71) Anthracene	11.56	178	1328755	49.108	nq	99
72) Carbazole	11.72	167	1214407	47.024	nq	100
73) Di-n-butylphthalate	12.02	149	1302685	42.348	nq	99
74) Fluoranthene	12.70	202	1306049	47.432	nq	99
76) Benzidine	12.84	184	280572	27.585	nq	98
77) Pyrene	12.93	202	1282786	49.997	nq	100
79) Butylbenzylphthalate	13.53	149	563277	46.599	nq	98
80) Benzo(a)anthracene	14.13	228	946147	42.366	nq	99
81) 3,3'-Dichlorobenzidine	14.09	252	181833	24.598	nq	98
82) Chrysene	14.17	228	1069205	48.880	nq	99
83) Bis(2-ethylhexyl)phthalate	14.07	149	728311	46.632	nq	100
84) Di-n-octyl phthalate	14.70	149	1184992	44.107	nq	# 94
85) Indeno(1,2,3-cd)pyrene	17.26	276	859995	37.940	nq	97
87) Benzo(b)fluoranthene	15.22	252	747473	39.673	nq	98
88) Benzo(k)fluoranthene	15.24	252	1007540	56.769	nq	99
89) Benzo(a)pyrene	15.61	252	827707	47.407	nq	99
90) Dibenzo(a,h)anthracene	17.27	278	703711	43.593	nq	98
91) Benzo(g,h,i)perylene	17.74	276	705468	43.631	nq	96

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

