

Data Path : U:\HPCHEM1\BNA F\DATA\BF020518\
 Data File : BF102743.D
 Acq On : 5 Feb 2018 17:38
 Operator : SJ/JU
 Sample : J1399-02MSD
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-01-020218-AMSD

Manual Integrations
 APPROVED

Sohil

Quant Time: Feb 06 11:20:01 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 05 12:44:16 2018
 Response via : Initial Calibration

2/6/2018 2:25:37 PM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	221497	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	902742	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	397583	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	611597	20.00	ng	0.00
75) Chrysene-d12	14.04	240	353385	20.00	ng	0.00
86) Perylene-d12	15.52	264	304199	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.52	112	1658194	113.69	ng	0.02
7) Phenol-d6	6.52	99	1832382	104.34	ng	0.01
23) Nitrobenzene-d5	7.45	82	1353796	104.03	ng	0.00
41) 2,4,6-Tribromophenol	10.72	330	429548	134.23	ng	0.00
44) 2-Fluorobiphenyl	9.25	172	2242721	93.60	ng	0.00
78) Terphenyl-d14	12.99	244	1596889	107.00	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.88	88	254649m	35.349	ng	
3) Pyridine	3.59	79	407714m	20.485	ng	
4) n-Nitrosodimethylamine	3.53	42	381720m	44.826	ng	
6) Aniline	6.56	93	160119	6.174	ng	96
8) 2-Chlorophenol	6.67	128	631351	42.047	ng	93
9) Benzaldehyde	6.43	77	222242	17.041	ng	97
10) Phenol	6.53	94	668565	35.524	ng	96
11) bis(2-Chloroethyl)ether	6.62	93	687003	43.869	ng	92
12) 1,3-Dichlorobenzene	6.82	146	678971	39.048	ng	99
13) 1,4-Dichlorobenzene	6.90	146	683312	39.450	ng	99
14) 1,2-Dichlorobenzene	7.05	146	628911	38.983	ng	99
15) Benzyl Alcohol	7.02	79	313616	23.228	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.15	45	1220664	41.032	ng	95
17) 2-Methylphenol	7.13	107	549391	39.666	ng	98
18) Hexachloroethane	7.39	117	236171	39.762	ng	92
19) n-Nitroso-di-n-propylamine	7.30	70	456591	39.434	ng	92
20) 3+4-Methylphenols	7.29	107	610262	35.730	ng	# 73
22) Acetophenone	7.29	105	895638	40.543	ng	# 94
24) Nitrobenzene	7.47	77	713817	46.635	ng	96
25) Isophorone	7.70	82	1332755	44.477	ng	100
26) 2-Nitrophenol	7.78	139	355020	54.927	ng	98
27) 2,4-Dimethylphenol	7.82	122	622719	49.189	ng	99
28) bis(2-Chloroethoxy)methane	7.91	93	620644	34.964	ng	99
29) 2,4-Dichlorophenol	8.02	162	558280	48.510	ng	97
30) 1,2,4-Trichlorobenzene	8.10	180	548916	42.162	ng	98
31) Naphthalene	8.19	128	1960729	44.455	ng	99
32) Benzoic acid	7.94	122	453730	49.940	ng	98
33) 4-Chloroaniline	8.23	127	56625	2.980	ng	96
34) Hexachlorobutadiene	8.30	225	278527	41.749	ng	99
35) Caprolactam	8.61	113	152745m	38.217	ng	
36) 4-Chloro-3-methylphenol	8.72	107	590704	45.682	ng	99
37) 2-Methylnaphthalene	8.88	142	1258004	43.564	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.05	216	476900	44.053	ng	99
40) Hexachlorocyclopentadiene	9.03	237	268485	52.171	ng	97

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42) 2,4,6-Trichlorophenol	9.16	196	360351	50.679	ng	99
43) 2,4,5-Trichlorophenol	9.20	196	384653	47.997	ng	99
45) 1,1'-Biphenyl	9.35	154	1435632	42.871	ng	98
46) 2-Chloronaphthalene	9.37	162	1126864	42.743	ng	98
47) 2-Nitroaniline	9.46	65	362042	44.756	ng	91
48) Acenaphthylene	9.79	152	1621645	39.603	ng	99
49) Dimethylphthalate	9.65	163	1376687	44.409	ng	100
50) 2,6-Dinitrotoluene	9.70	165	300615	50.970	ng	# 91
51) Acenaphthene	9.96	154	1086354	44.363	ng	98
52) 3-Nitroaniline	9.87	138	96668	13.321	ng	98
53) 2,4-Dinitrophenol	9.98	184	167420	82.533	ng	# 22
54) Dibenzofuran	10.13	168	1502984	43.553	ng	99
55) 4-Nitrophenol	10.03	139	431650	72.763	ng	94
56) 2,4-Dinitrotoluene	10.11	165	388830	49.083	ng	97
57) Fluorene	10.47	166	1077933	42.931	ng	100
58) 2,3,4,6-Tetrachlorophenol	10.25	232	277565	49.971	ng	99
59) Diethylphthalate	10.35	149	1309076	42.766	ng	99
60) 4-Chlorophenyl-phenylether	10.46	204	494133	44.488	ng	99
61) 4-Nitroaniline	10.49	138	158912	23.005	ng	91
62) Azobenzene	10.62	77	845912	27.663	ng	95
64) 4,6-Dinitro-2-methylphenol	10.52	198	115731	46.147	ng	82
65) n-Nitrosodiphenylamine	10.58	169	393587	18.245	ng	100
66) 4-Bromophenyl-phenylether	10.95	248	315344	48.743	ng	97
67) Hexachlorobenzene	11.02	284	318329	44.602	ng	# 88
68) Atrazine	11.10	200	22952	3.606	ng	98
69) Pentachlorophenol	11.21	266	349469	83.413	ng	98
70) Phenanthrene	11.43	178	1471520	43.201	ng	99
71) Anthracene	11.49	178	1495660	43.172	ng	100
72) Carbazole	11.63	167	498617	14.691	ng	99
73) Di-n-butylphthalate	11.97	149	1900163	46.088	ng	99
74) Fluoranthene	12.62	202	1290693	37.662	ng	99
77) Pyrene	12.85	202	1230780	44.415	ng	100
79) Butylbenzylphthalate	13.46	149	688447	51.854	ng	98
80) Benzo(a)anthracene	14.03	228	1012730	45.568	ng	99
81) 3,3'-Dichlorobenzidine	13.99	252	28522	3.461	ng	99
82) Chrysene	14.07	228	981468	43.809	ng	99
83) Bis(2-ethylhexyl)phthalate	14.02	149	920547	54.228	ng	99
84) Di-n-octyl phthalate	14.63	149	1523663	59.777	ng	98
85) Indeno(1,2,3-cd)pyrene	17.02	276	865547	46.582	ng	98
87) Benzo(b)fluoranthene	15.09	252	1016265	51.529	ng	98
88) Benzo(k)fluoranthene	15.12	252	954052	50.628	ng	100
89) Benzo(a)pyrene	15.46	252	901361	50.774	ng	98
90) Dibenzo(a,h)anthracene	17.04	278	753597	52.300	ng	99
91) Benzo(g,h,i)perylene	17.47	276	690678	48.972	ng	97

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

