

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122217\
 Data File : BF101662.D
 Acq On : 22 Dec 2017 19:15
 Operator : SJ/JU
 Sample : I7002-01MS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EPE-106-3DS(16-20)MS

Quant Time: Dec 22 22:34:40 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 22 16:07:58 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	142330	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	541651	20.00	ng	0.00
38) Acenaphthene-d10	9.83	164	227317	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	373981	20.00	ng	0.00
75) Chrysene-d12	13.97	240	269012	20.00	ng	0.00
86) Perylene-d12	15.43	264	258581	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.42	112	1123383	117.57	ng	0.01
7) Phenol-d6	6.44	99	1293748	108.79	ng	0.00
23) Nitrobenzene-d5	7.36	82	816690	83.93	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	247848	113.15	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	1176312	80.27	ng	0.00
78) Terphenyl-d14	12.90	244	917431	82.22	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.56	88	161888	32.973	ng	93
3) Pyridine	3.33	79	397999	29.176	ng	95
4) n-Nitrosodimethylamine	3.28	42	215332	34.284	ng	# 99
6) Aniline	6.46	93	241647	14.623	ng	# 80
8) 2-Chlorophenol	6.58	128	354171	35.236	ng	99
9) Benzaldehyde	6.35	77	193612	22.046	ng	96
10) Phenol	6.45	94	494562	36.489	ng	95
11) bis(2-Chloroethyl)ether	6.53	93	377874	35.543	ng	94
12) 1,3-Dichlorobenzene	6.73	146	392105	34.565	ng	100
13) 1,4-Dichlorobenzene	6.80	146	402014	34.703	ng	98
14) 1,2-Dichlorobenzene	6.96	146	358782	34.408	ng	99
15) Benzyl Alcohol	6.94	79	267461	32.677	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.06	45	613671	37.716	ng	71
17) 2-Methylphenol	7.06	107	295481	31.959	ng	# 87
18) Hexachloroethane	7.29	117	123702	30.713	ng	94
19) n-Nitroso-di-n-propylamine	7.20	70	255937	32.773	ng	94
20) 3+4-Methylphenols	7.22	107	401434	40.973	ng	# 87
22) Acetophenone	7.20	105	511913	41.420	ng	# 95
24) Nitrobenzene	7.38	77	400414	37.182	ng	100
25) Isophorone	7.62	82	723532	37.066	ng	97
26) 2-Nitrophenol	7.69	139	189312	40.918	ng	96
27) 2,4-Dimethylphenol	7.73	122	318056	40.465	ng	97
28) bis(2-Chloroethoxy)methane	7.82	93	456070	36.782	ng	99
29) 2,4-Dichlorophenol	7.95	162	301055	38.769	ng	98
30) 1,2,4-Trichlorobenzene	8.02	180	290643	35.467	ng	94
31) Naphthalene	8.09	128	978583	35.887	ng	100
32) Benzoic acid	7.85	122	14559	10.254	ng	96
33) 4-Chloroaniline	8.16	127	125609	10.598	ng	99
34) Hexachlorobutadiene	8.20	225	169861	35.839	ng	96
35) Caprolactam	8.53	113	98293	36.454	ng	95
36) 4-Chloro-3-methylphenol	8.65	107	302843	35.483	ng	99
37) 2-Methylnaphthalene	8.79	142	624902	35.174	ng	100
39) 1,2,4,5-Tetrachlorobenzene	8.95	216	295865	38.550	ng	97
40) Hexachlorocyclopentadiene	8.93	237	2094	13.193	ng	86

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42) 2,4,6-Trichlorophenol	9.07	196	186076	40.272	ng	100
43) 2,4,5-Trichlorophenol	9.12	196	185201	36.608	ng	98
45) 1,1'-Biphenyl	9.25	154	750050	37.206	ng	99
46) 2-Chloronaphthalene	9.27	162	557310	36.130	ng	98
47) 2-Nitroaniline	9.39	65	205703	38.603	ng	94
48) Acenaphthylene	9.69	152	803567	32.982	ng	100
49) Dimethylphthalate	9.55	163	745288	40.761	ng	99
50) 2,6-Dinitrotoluene	9.62	165	133820	36.010	ng	# 81
51) Acenaphthene	9.86	154	490806	33.530	ng	98
52) 3-Nitroaniline	9.80	138	103987	22.444	ng	98
53) 2,4-Dinitrophenol	9.95	184	7096	15.095	ng	# 21
54) Dibenzofuran	10.03	168	691446	37.170	ng	99
55) 4-Nitrophenol	9.98	139	142897	70.732	ng	95
56) 2,4-Dinitrotoluene	10.04	165	165353	39.493	ng	# 97
57) Fluorene	10.37	166	552188	35.090	ng	100
58) 2,3,4,6-Tetrachlorophenol	10.16	232	145416	40.007	ng	# 86
59) Diethylphthalate	10.24	149	624697	34.500	ng	98
60) 4-Chlorophenyl-phenylether	10.36	204	271234	35.964	ng	94
61) 4-Nitroaniline	10.42	138	127808	27.444	ng	98
62) Azobenzene	10.53	77	618633	32.981	ng	96
64) 4,6-Dinitro-2-methylphenol	10.46	198	22162	11.613	ng	92
65) n-Nitrosodiphenylamine	10.49	169	514534	37.261	ng	95
66) 4-Bromophenyl-phenylether	10.85	248	163530	38.916	ng	93
67) Hexachlorobenzene	10.93	284	158553	35.650	ng	92
68) Atrazine	11.02	200	159489	38.734	ng	96
69) Pentachlorophenol	11.13	266	123956	72.354	ng	98
70) Phenanthrene	11.34	178	742383	35.696	ng	99
71) Anthracene	11.39	178	748987	35.256	ng	100
72) Carbazole	11.56	167	677903	34.082	ng	99
73) Di-n-butylphthalate	11.87	149	898509	37.219	ng	99
74) Fluoranthene	12.53	202	715605	32.781	ng	99
76) Benzidine	12.66	184	323546	28.477	ng	99
77) Pyrene	12.76	202	723438	35.833	ng	100
79) Butylbenzylphthalate	13.37	149	334000	36.562	ng	95
80) Benzo(a)anthracene	13.96	228	629289	35.426	ng	98
81) 3,3'-Dichlorobenzidine	13.92	252	172877	29.847	ng	# 98
82) Chrysene	13.99	228	599462	35.742	ng	100
83) Bis(2-ethylhexyl)phthalate	13.92	149	446232	38.565	ng	99
84) Di-n-octyl phthalate	14.53	149	762423	36.947	ng	97
85) Indeno(1,2,3-cd)pyrene	16.92	276	524066	35.089	ng	96
87) Benzo(b)fluoranthene	15.01	252	570625	36.205	ng	98
88) Benzo(k)fluoranthene	15.04	252	512781	33.462	ng	98
89) Benzo(a)pyrene	15.37	252	520246	35.806	ng	99
90) Dibenzo(a,h)anthracene	16.91	278	445011	35.673	ng	98
91) Benzo(g,h,i)perylene	17.36	276	433962	35.131	ng	95

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

