

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060816\
 Data File : BF087839.D
 Acq On : 9 Jun 2016 7:43
 Operator : UM/SJ
 Sample : H3405-02MSD
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-13-COMP-0.5-11.0MSD

Manual Integrations
 APPROVED

sohil
 6/9/2016 7:58:50 PM

Quant Time: Jun 09 08:36:07 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 07 18:51:55 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	110042	20.00	ng	-0.01
21) Naphthalene-d8	8.11	136	428132	20.00	ng	-0.01
38) Acenaphthene-d10	9.87	164	180015	20.00	ng	-0.01
63) Phenanthrene-d10	11.36	188	276670	20.00	ng	0.00
75) Chrysene-d12	13.99	240	194547	20.00	ng	0.00
86) Perylene-d12	15.42	264	173119	20.00	ng	0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.43	112	942852	146.48	ng	0.00
7) Phenol-d6	6.47	99	1145767	146.87	ng	-0.01
23) Nitrobenzene-d5	7.39	82	670847	91.50	ng	-0.02
41) 2,4,6-Tribromophenol	10.66	330	207495	109.16	ng	-0.01
44) 2-Fluorobiphenyl	9.20	172	1091018	95.92	ng	-0.01
78) Terphenyl-d14	12.95	244	849017	99.31	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.46	88	106251	33.97	ng	# 33
3) Pyridine	3.19	79	314468	42.58	ng	96
4) n-Nitrosodimethylamine	3.11	42	146857	46.96	ng	94
6) Aniline	6.49	93	257577	23.20	ng	# 55
8) 2-Chlorophenol	6.62	128	346256	45.45	ng	90
9) Benzaldehyde	6.36	77	21015	3.58	ng	86
10) Phenol	6.48	94	300471	36.85	ng	# 69
11) bis(2-Chloroethyl)ether	6.57	93	312346	45.66	ng	88
12) 1,3-Dichlorobenzene	6.76	146	319796	39.12	ng	99
13) 1,4-Dichlorobenzene	6.84	146	339234	41.20	ng	99
14) 1,2-Dichlorobenzene	7.00	146	314489	41.99	ng	100
15) Benzyl Alcohol	6.97	79	238735	39.12	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.11	45	399030	43.18	ng	96
17) 2-Methylphenol	7.08	107	252590	40.88	ng	100
18) Hexachloroethane	7.34	117	100901	34.53	ng	# 81
19) n-Nitroso-di-n-propylamine	7.26	70	215505	43.18	ng	# 80
20) 3+4-Methylphenols	7.24	107	304593	42.25	ng	# 86
22) Acetophenone	7.24	105	407112	42.55	ng	# 90
24) Nitrobenzene	7.42	77	353520	49.78	ng	95
25) Isophorone	7.66	82	595414	43.84	ng	97
26) 2-Nitrophenol	7.72	139	150857	39.65	ng	97
27) 2,4-Dimethylphenol	7.77	122	254928	36.39	ng	94
28) bis(2-Chloroethoxy)methane	7.87	93	376631	44.71	ng	98
29) 2,4-Dichlorophenol	7.98	162	276320	47.25	ng	93
30) 1,2,4-Trichlorobenzene	8.06	180	264000	41.46	ng	99
31) Naphthalene	8.14	128	877153	41.40	ng	100
32) Benzoic acid	7.84	122	24258m	6.29	ng	
33) 4-Chloroaniline	8.18	127	150702	15.93	ng	98
34) Hexachlorobutadiene	8.25	225	135296	40.28	ng	97
35) Caprolactam	8.56	113	71953	37.14	ng	95
36) 4-Chloro-3-methylphenol	8.67	107	240757	38.49	ng	98
37) 2-Methylnaphthalene	8.83	142	568103	40.68	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.99	216	230407	51.84	ng	# 1
40) Hexachlorocyclopentadiene	8.98	237	37496	12.91	ng	98

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42) 2,4,6-Trichlorophenol	9.11	196	161130	44.76	nq	97
43) 2,4,5-Trichlorophenol	9.15	196	150378	40.15	nq #	85
45) 1,1'-Biphenyl	9.30	154	647471	48.39	nq	95
46) 2-Chloronaphthalene	9.32	162	492986	42.32	nq #	91
47) 2-Nitroaniline	9.42	65	161214	46.91	nq	88
48) Acenaphthylene	9.74	152	792656	42.78	nq	99
49) Dimethylphthalate	9.60	163	942750	72.30	nq	99
50) 2,6-Dinitrotoluene	9.66	165	119736	40.09	nq #	52
51) Acenaphthene	9.91	154	459578	39.76	nq	100
52) 3-Nitroaniline	9.83	138	124510	36.13	nq	87
53) 2,4-Dinitrophenol	9.93	184	12908	12.42	nq #	85
54) Dibenzofuran	10.08	168	674715	42.38	nq	94
55) 4-Nitrophenol	9.99	139	189984	71.03	nq	95
56) 2,4-Dinitrotoluene	10.06	165	150408	39.19	nq #	56
57) Fluorene	10.42	166	497558	46.40	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.19	232	111496	37.88	nq #	54
59) Diethylphthalate	10.31	149	594763	45.06	nq	97
60) 4-Chlorophenyl-phenylether	10.41	204	206319	38.97	nq	93
61) 4-Nitroaniline	10.44	138	130089	39.80	nq	90
62) Azobenzene	10.57	77	511524	39.19	nq	96
64) 4,6-Dinitro-2-methylphenol	10.47	198	18716	12.51	nq	82
65) n-Nitrosodiphenylamine	10.54	169	457420	49.83	nq	100
66) 4-Bromophenyl-phenylether	10.90	248	125083	42.14	nq #	84
67) Hexachlorobenzene	10.97	284	143789	46.62	nq #	76
68) Atrazine	11.06	200	110691	39.82	nq	92
69) Pentachlorophenol	11.16	266	149240	91.80	nq	100
70) Phenanthrene	11.38	178	666664	45.92	nq	99
71) Anthracene	11.43	178	593883	40.55	nq	99
72) Carbazole	11.59	167	626628	45.73	nq	99
73) Di-n-butylphthalate	11.93	149	788962	45.48	nq #	97
74) Fluoranthene	12.56	202	625979	40.86	nq	96
76) Benzidine	12.69	184	150796	27.99	nq	99
77) Pyrene	12.79	202	622045	43.09	nq	99
79) Butylbenzylphthalate	13.42	149	306689	48.05	nq #	81
80) Benzo(a)anthracene	13.98	228	535452	48.30	nq	99
81) 3,3'-Dichlorobenzidine	13.94	252	143410	48.10	nq #	95
82) Chrysene	14.02	228	545256	52.28	nq	98
83) Bis(2-ethylhexyl)phthalate	13.99	149	420401	54.11	nq #	97
84) Di-n-octyl phthalate	14.59	149	727339	57.61	nq #	100
85) Indeno(1,2,3-cd)pyrene	16.80	276	413661	42.69	nq #	100
87) Benzo(b)fluoranthene	15.01	252	450286m	37.32	nq	
88) Benzo(k)fluoranthene	15.04	252	460655m	50.61	nq	
89) Benzo(a)pyrene	15.36	252	413390	42.35	nq #	94
90) Dibenzo(a,h)anthracene	16.82	278	345960	38.16	nq	99
91) Benzo(g,h,i)perylene	17.22	276	339338	36.38	nq	98

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

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