

Data Path : V:\HPCHEM1\BNA F\DATA\BF070916\
 Data File : BF088866.D
 Acq On : 9 Jul 2016 12:38
 Operator : UM/SJ
 Sample : PB91808BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB91808BS

Manual Integrations

sohil
 7/11/2016 7:15:33 PM

Quant Time: Jul 11 05:27:23 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 08 18:51:20 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	70015	20.00	ng	-0.01
21) Naphthalene-d8	8.00	136	259866	20.00	ng	-0.02
38) Acenaphthene-d10	9.75	164	125556	20.00	ng	-0.02
63) Phenanthrene-d10	11.22	188	257449	20.00	ng	-0.02
75) Chrysene-d12	13.85	240	230937	20.00	ng	-0.02
86) Perylene-d12	15.22	264	201344	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.30	112	63378	13.57	ng	-0.02
7) Phenol-d6	6.34	99	76831	12.73	ng	-0.03
23) Nitrobenzene-d5	7.28	82	50916	10.27	ng	-0.02
41) 2,4,6-Tribromophenol	10.54	330	20460	17.55	ng	-0.02
44) 2-Fluorobiphenyl	9.07	172	91441	11.50	ng	-0.02
78) Terphenyl-d14	12.81	244	101027	9.74	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.27	88	9182	3.96	ng	# 34
8) 2-Chlorophenol	6.50	128	10620	2.12	ng	83
12) 1,3-Dichlorobenzene	6.66	146	11073m	2.00	ng	
13) 1,4-Dichlorobenzene	6.73	146	11237m	2.05	ng	
14) 1,2-Dichlorobenzene	6.89	146	11977	2.33	ng	97
17) 2-Methylphenol	6.98	107	8542	2.09	ng	97
18) Hexachloroethane	7.23	117	4486	2.11	ng	90
19) n-Nitroso-di-n-propylamine	7.13	70	8441	2.08	ng	# 87
20) 3+4-Methylphenols	7.13	107	11488	2.34	ng	# 83
22) Acetophenone	7.13	105	16107	2.55	ng	# 91
24) Nitrobenzene	7.29	77	11149	2.23	ng	# 79
25) Isophorone	7.54	82	21862	2.38	ng	99
26) 2-Nitrophenol	7.62	139	4772	2.33	ng	# 85
27) 2,4-Dimethylphenol	7.67	122	9820	2.30	ng	97
28) bis(2-Chloroethoxy)methane	7.76	93	14669	2.45	ng	# 96
29) 2,4-Dichlorophenol	7.86	162	8545	2.48	ng	100
30) 1,2,4-Trichlorobenzene	7.94	180	9578	2.53	ng	95
31) Naphthalene	8.02	128	33107	2.58	ng	99
34) Hexachlorobutadiene	8.15	225	5197	2.56	ng	99
35) Caprolactam	8.39	113	2564	2.19	ng	90
36) 4-Chloro-3-methylphenol	8.56	107	10188	2.50	ng	98
37) 2-Methylnaphthalene	8.72	142	20711	2.51	ng	96
39) 1,2,4,5-Tetrachlorobenzene	8.88	216	8963	2.15	ng	# 1
40) Hexachlorocyclopentadiene	8.87	237	8585	4.19	ng	95
42) 2,4,6-Trichlorophenol	8.99	196	6532	2.37	ng	97
43) 2,4,5-Trichlorophenol	9.03	196	5527	2.33	ng	# 83
45) 1,1'-Biphenyl	9.18	154	27372	2.63	ng	99
46) 2-Chloronaphthalene	9.20	162	21102	2.59	ng	99
47) 2-Nitroaniline	9.29	65	5991	2.07	ng	93
48) Acenaphthylene	9.61	152	34633	2.67	ng	98
49) Dimethylphthalate	9.47	163	24128	2.70	ng	99
50) 2,6-Dinitrotoluene	9.53	165	4613	2.33	ng	# 33
51) Acenaphthene	9.78	154	23430	2.94	ng	98
52) 3-Nitroaniline	9.70	138	5812	2.31	ng	95

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
53) 2,4-Dinitrophenol	9.80	184	3207	3.66	nq	# 78
54) Dibenzofuran	9.95	168	31365	3.01	nq	92
55) 4-Nitrophenol	9.86	139	11754	6.14	nq	97
56) 2,4-Dinitrotoluene	9.93	165	5959	2.30	nq	# 17
57) Fluorene	10.30	166	25333	3.15	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.08	232	4591	2.50	nq	# 54
59) Diethylphthalate	10.17	149	22967	2.56	nq	96
60) 4-Chlorophenyl-phenylether	10.30	204	10942	2.93	nq	94
61) 4-Nitroaniline	10.30	138	5060	2.09	nq	76
62) Azobenzene	10.44	77	26190	2.56	nq	90
65) n-Nitrosodiphenylamine	10.41	169	21293	2.44	nq	99
66) 4-Bromophenyl-phenylether	10.78	248	5877	2.27	nq	# 65
67) Hexachlorobenzene	10.84	284	7180	2.50	nq	# 80
68) Atrazine	10.94	200	6726	2.48	nq	95
69) Pentachlorophenol	11.04	266	7220	4.25	nq	93
70) Phenanthrene	11.24	178	35615	2.53	nq	98
71) Anthracene	11.30	178	37879	2.71	nq	98
72) Carbazole	11.46	167	30274	2.30	nq	98
73) Di-n-butylphthalate	11.81	149	37124	2.42	nq	97
74) Fluoranthene	12.43	202	35476	2.49	nq	95
76) Benzidine	12.56	184	25729	2.20	nq	99
80) Benzo(a)anthracene	13.84	228	33482	2.25	nq	98
82) Chrysene	13.87	228	30640	2.08	nq	97
83) Bis(2-ethylhexyl)phthalate	13.85	149	20898	2.10	nq	99
85) Indeno(1,2,3-cd)pyrene	16.51	276	23833	2.11	nq	# 100
87) Benzo(b)fluoranthene	14.85	252	30200m	2.39	nq	
88) Benzo(k)fluoranthene	14.87	252	22980m	2.09	nq	
89) Benzo(a)pyrene	15.17	252	24558	2.30	nq	96
90) Dibenzo(a,h)anthracene	16.53	278	19711	2.21	nq	98
91) Benzo(g,h,i)perylene	16.89	276	19241	2.08	nq	100

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

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