

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042517\
 Data File : BF094611.D
 Acq On : 25 Apr 2017 18:00
 Operator : SJ/MA
 Sample : I2835-01MS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MARJON-SEMS

Manual Integrations
 APPROVED

mohammad
 4/26/2017 2:21:01 PM

Quant Time: Apr 26 03:10:57 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF041717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 17 14:59:23 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.75	152	126544	20.00	ng	0.00
21) Naphthalene-d8	7.25	136	492384	20.00	ng	-0.01
38) Acenaphthene-d10	9.38	164	214503	20.00	ng	-0.01
63) Phenanthrene-d10	11.20	188	341419	20.00	ng	-0.02
75) Chrysene-d12	14.46	240	249808	20.00	ng	0.00
86) Perylene-d12	16.08	264	235417	20.00	ng	0.00

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System Monitoring Compounds

5) 2-Fluorophenol	4.31	112	853673	111.92	ng	0.00
7) Phenol-d6	5.39	99	1022364	113.70	ng	0.00
23) Nitrobenzene-d5	6.41	82	637139	79.90	ng	0.00
41) 2,4,6-Tribromophenol	10.36	330	179761	107.14	ng	-0.01
44) 2-Fluorobiphenyl	8.57	172	1158220	79.45	ng	-0.01
78) Terphenyl-d14	13.18	244	786700	84.18	ng	-0.01

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Target Compounds

						Qvalue
2) 1,4-Dioxane	2.10	88	112786	25.43	ng	97
3) Pyridine	2.64	79	287242	29.63	ng	99
4) n-Nitrosodimethylamine	2.57	42	138852	34.61	ng	90
6) Aniline	5.38	93	372322	31.16	ng	# 90
8) 2-Chlorophenol	5.53	128	336493	38.77	ng	94
9) Benzaldehyde	5.25	77	102058	14.92	ng	99
10) Phenol	5.40	94	467042	42.28	ng	92
11) bis(2-Chloroethyl)ether	5.47	93	309215	38.32	ng	95
12) 1,3-Dichlorobenzene	5.69	146	355541	36.97	ng	99
13) 1,4-Dichlorobenzene	5.78	146	361157	37.33	ng	98
14) 1,2-Dichlorobenzene	5.95	146	336868	37.80	ng	95
15) Benzyl Alcohol	5.94	79	253282	37.97	ng	99
16) 2,2'-oxybis(1-Chloropropan	6.09	45	406730	38.44	ng	73
17) 2-Methylphenol	6.08	107	265368	38.82	ng	92
18) Hexachloroethane	6.34	117	114195	34.43	ng	88
19) n-Nitroso-di-n-propylamine	6.25	70	232979	39.10	ng	95
20) 3+4-Methylphenols	6.26	107	366892	39.50	ng	# 63
22) Acetophenone	6.23	105	471593	39.39	ng	# 85
24) Nitrobenzene	6.43	77	331241	39.45	ng	91
25) Isophorone	6.72	82	591858	40.04	ng	95
26) 2-Nitrophenol	6.80	139	166013	36.80	ng	98
27) 2,4-Dimethylphenol	6.88	122	287678	39.26	ng	99
28) bis(2-Chloroethoxy)methane	6.98	93	383788	40.14	ng	99
29) 2,4-Dichlorophenol	7.10	162	272021	39.90	ng	98
30) 1,2,4-Trichlorobenzene	7.19	180	277442	39.37	ng	98
31) Naphthalene	7.28	128	935309	39.51	ng	99
33) 4-Chloroaniline	7.35	127	177636	18.04	ng	# 94
34) Hexachlorobutadiene	7.43	225	137433	39.11	ng	97
35) Caprolactam	7.80	113	70478m	32.91	ng	
36) 4-Chloro-3-methylphenol	7.98	107	273755	38.32	ng	88
37) 2-Methylnaphthalene	8.12	142	632832	39.08	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.32	216	249187	40.80	ng	99
40) Hexachlorocyclopentadiene	8.31	237	73492	29.74	ng	98
42) 2,4,6-Trichlorophenol	8.48	196	169147	39.43	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	
43) 2,4,5-Trichlorophenol	8.54	196	169017	38.37	nq	#	92
45) 1,1'-Biphenyl	8.69	154	715001	40.80	nq		98
46) 2-Chloronaphthalene	8.71	162	559643	40.16	nq		94
47) 2-Nitroaniline	8.85	65	164164	40.95	nq	#	79
48) Acenaphthylene	9.21	152	841382	38.73	nq		98
49) Dimethylphthalate	9.08	163	790477	49.45	nq		99 mohammad
50) 2,6-Dinitrotoluene	9.15	165	142230	39.64	nq		98
51) Acenaphthene	9.42	154	509357	39.15	nq		99
52) 3-Nitroaniline	9.35	138	128882	31.05	nq		92
53) 2,4-Dinitrophenol	9.51	184	3643	7.49	nq	#	74
54) Dibenzofuran	9.64	168	747699	39.54	nq		97 4/26/2017 2:21:01 PM
55) 4-Nitrophenol	9.61	139	154331m	52.62	nq		
56) 2,4-Dinitrotoluene	9.65	165	179630	40.80	nq	#	83
57) Fluorene	10.06	166	567830	38.56	nq		100
58) 2,3,4,6-Tetrachlorophenol	9.81	232	114548	36.28	nq		93
59) Diethylphthalate	9.95	149	623683	41.35	nq		99
60) 4-Chlorophenyl-phenylether	10.07	204	250811	40.93	nq		94
61) 4-Nitroaniline	10.11	138	141408	33.51	nq		95
62) Azobenzene	10.26	77	587945	40.15	nq		97
64) 4,6-Dinitro-2-methylphenol	10.15	198	10948	4.89	nq	#	68
65) n-Nitrosodiphenylamine	10.22	169	523456	44.08	nq		99
66) 4-Bromophenyl-phenylether	10.67	248	149584	44.12	nq		96
67) Hexachlorobenzene	10.74	284	139115	40.72	nq	#	85
68) Atrazine	10.89	200	152123	42.82	nq		95
69) Pentachlorophenol	10.99	266	99844	73.60	nq		97
70) Phenanthrene	11.24	178	765785	39.83	nq		98
71) Anthracene	11.30	178	783863	39.41	nq		99
72) Carbazole	11.51	167	718915	38.51	nq		98
73) Di-n-butylphthalate	11.96	149	933422	45.42	nq		98
74) Fluoranthene	12.69	202	690373	34.65	nq		99
76) Benzidine	12.88	184	350880	34.69	nq	#	93
77) Pyrene	12.97	202	681206	39.03	nq		98
79) Butylbenzylphthalate	13.79	149	326842	42.73	nq		89
80) Benzo(a)anthracene	14.45	228	561435	38.67	nq		98
81) 3,3'-Dichlorobenzidine	14.42	252	191221	37.20	nq	#	96
82) Chrysene	14.49	228	520686	39.24	nq		99
83) Bis(2-ethylhexyl)phthalate	14.51	149	458487	44.65	nq	#	99
84) Di-n-octyl phthalate	15.27	149	754883	43.82	nq		98
85) Indeno(1,2,3-cd)pyrene	17.36	276	489450	38.77	nq		100
87) Benzo(b)fluoranthene	15.68	252	542197	37.65	nq		98
88) Benzo(k)fluoranthene	15.71	252	546436	40.10	nq		96
89) Benzo(a)pyrene	16.03	252	523842	39.10	nq		99
90) Dibenzo(a,h)anthracene	17.37	278	414570	37.27	nq		96
91) Benzo(g,h,i)perylene	17.73	276	382676	33.79	nq		99

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

