

Data Path : Z:\HPCHEM1\BNA M\DATA\BM092717\
 Data File : BM011750.D
 Acq On : 28 Sep 2017 07:23
 Operator : SJ/JU
 Sample : I5377-23
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CB3J3

Manual Integrations
 APPROVED

Sohil
 9/29/2017 2:23:15 PM

Quant Time: Sep 28 08:04:05 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM092717.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 28 05:09:12 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	295409	20.00	ng/ul	0.00
18) Naphthalene-d8	10.74	136	1308127	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.57	164	730060	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.32	188	1561918	20.00	ng/ul	0.00
75) Chrysene-d12	21.48	240	1291751	20.00	ng/ul	0.00
83) Perylene-d12	23.84	264	1104424	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.36	96	9723	1.47	ng/uL	0.00
5) Phenol-d5	7.09	99	394089	13.56	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.26	67	277486	13.17	ng/ul	0.00
9) 2-Chlorophenol-d4	7.46	132	298350	13.80	ng/ul	0.00
13) 4-Methylphenol-d8	8.63	113	350303	15.39	ng/ul	0.00
19) Nitrobenzene-d5	9.10	128	159729	15.65	ng/ul	0.00
22) 2-Nitrophenol-d4	9.82	143	178884	16.44	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.36	165	351734	16.93	ng/ul	0.00
29) 4-Chloroaniline-d4	10.87	131	263479	11.47	ng/ul	0.00
43) Dimethylphthalate-d6	13.97	166	1179455	18.84	ng/ul	0.00
46) Acenaphthylene-d8	14.26	160	1399503	18.48	ng/ul	0.00
51) 4-Nitrophenol-d4	14.76	143	151104m	14.12	ng/ul	0.00
57) Fluorene-d10	15.56	176	1013566	18.80	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.68	200	135293	12.79	ng/ul	0.00
70) Anthracene-d10	17.41	188	1534333	19.30	ng/ul	0.00
76) Pyrene-d10	19.70	212	1624680	26.35	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.69	264	1035615	19.58	ng/ul	0.00

Target Compounds

					Ovalue	
28) Naphthalene	10.79	128	131708	1.926	ng/ul	99
34) 2-Methylnaphthalene	12.40	142	60738	1.253	ng/ul	94
44) Dimethylphthalate	14.02	163	236685	3.952	ng/ul	99
49) Acenaphthene	14.63	153	97015	1.875	ng/ul	98
53) Dibenzofuran	14.97	168	91977	1.295	ng/ul	90
58) Fluorene	15.62	166	104562	1.738	ng/ul#	95
69) Phenanthrene	17.36	178	1924148	22.011	ng/ul	98
71) Anthracene	17.44	178	378893	4.207	ng/ul	97
72) Carbazole	17.72	167	138037	1.749	ng/ul	98
74) Fluoranthene	19.37	202	2907830	27.244	ng/ul#	91
77) Pyrene	19.73	202	2758844	35.839	ng/ul#	88
80) Benzo(a)anthracene	21.47	228	1310440	18.133	ng/ul	95
81) Bis(2-ethylhexyl)phthalate	21.37	149	57509	1.197	ng/ul#	97
82) Chrysene	21.52	228	1331033	19.528	ng/ul	96
85) Benzo(b)fluoranthene	23.13	252	1653095	25.369	ng/ul#	93
86) Benzo(k)fluoranthene	23.17	252	617912m	9.295	ng/ul	
88) Benzo(a)pyrene	23.74	252	1195080	18.989	ng/ul#	95
89) Indeno(1,2,3-cd)pyrene	26.27	276	892024	13.613	ng/ul#	93
90) Dibenzo(a,h)anthracene	26.27	278	265417	4.836	ng/ul#	85
91) Benzo(g,h,i)perylene	27.01	276	927130m	17.046	ng/ul	

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

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