

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061217\
 Data File : BM010552.D
 Acq On : 13 Jun 2017 01:21
 Operator : SJ/MA
 Sample : I3521-10DL 30X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F5C58DL

Manual Integrations
 APPROVED

mohammad
 6/13/2017 3:54:51 PM

Quant Time: Jun 13 04:52:00 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 13 03:33:28 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	102777	20.00	ng/ul	0.00
18) Naphthalene-d8	10.69	136	372428	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.52	164	255290	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.27	188	694204	20.00	ng/ul	0.00
75) Chrysene-d12	21.44	240	1030648	20.00	ng/ul	0.00
83) Perylene-d12	23.77	264	1044544	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
5) Phenol-d5	7.08	99	2709	0.35	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	7.23	67	1980m	0.45	ng/ul	0.00
9) 2-Chlorophenol-d4	7.43	132	3251m	0.52	ng/ul	0.00
13) 4-Methylphenol-d8	0.00	113	0d	0.00	ng/ul	
19) Nitrobenzene-d5	9.08	128	725	0.27	ng/ul	0.00
22) 2-Nitrophenol-d4	9.79	143	2353	0.74	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.33	165	3327	0.51	ng/ul	0.02
29) 4-Chloroaniline-d4	10.87	131	6469m	1.03	ng/ul	0.01
43) Dimethylphthalate-d6	13.94	166	17177	0.81	ng/ul	0.00
46) Acenaphthylene-d8	14.23	160	16127	0.65	ng/ul	0.00
51) 4-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul	
57) Fluorene-d10	15.52	176	15551	0.83	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.66	200	2012	0.41	ng/ul	0.01
70) Anthracene-d10	17.37	188	21533	0.64	ng/ul	0.00
76) Pyrene-d10	19.66	212	27478	0.64	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.61	264	32739	0.67	ng/ul	-0.02

Target Compounds

					Qvalue
28) Naphthalene	10.74	128	1618748	86.657 ng/ul	99
34) 2-Methylnaphthalene	12.34	142	558598	39.441 ng/ul	98
40) 1,1'-Biphenyl	13.36	154	161292	7.788 ng/ul	94
49) Acenaphthene	14.59	153	475080	28.014 ng/ul	97
53) Dibenzofuran	14.92	168	641223	25.571 ng/ul	99
58) Fluorene	15.57	166	630348	30.645 ng/ul	99
69) Phenanthrene	17.32	178	3117038	84.107 ng/ul	99
71) Anthracene	17.40	178	307083	7.937 ng/ul	96
72) Carbazole	17.69	167	139123	4.248 ng/ul	97
74) Fluoranthene	19.32	202	1837249	40.314 ng/ul#	96
77) Pyrene	19.69	202	1185534	22.322 ng/ul#	93
80) Benzo(a)anthracene	21.42	228	383801	6.604 ng/ul	99
82) Chrysene	21.47	228	302891	5.765 ng/ul	96
85) Benzo(b)fluoranthene	23.06	252	220515	3.628 ng/ul	100
86) Benzo(k)fluoranthene	23.10	252	67415m	1.161 ng/ul	
88) Benzo(a)pyrene	23.67	252	121219	2.012 ng/ul	96

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

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