

Data Path : Z:\HPCHEM1\BNA M\DATA\BM092217\
 Data File : BM011692.D
 Acq On : 22 Sep 2017 14:54
 Operator : SJ/JU
 Sample : I5283-01RE
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleID :
 CB3A6RE

Manual Integrations
 APPROVED

Sohil
 9/25/2017 6:23:31 PM

Quant Time: Sep 23 01:01:17 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM090817MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 22 23:48:55 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.96	152	379535	20.00	ng/ul	0.00
18) Naphthalene-d8	10.77	136	1718240	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.59	164	941868	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.33	188	1987347	20.00	ng/ul	0.00
78) Chrysene-d12	21.50	240	1463027	20.00	ng/ul	0.00
86) Perylene-d12	23.87	264	1287469	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.39	96	1904	0.23	ng/uL	0.00
5) Phenol-d5	7.12	99	19870	0.55	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.29	67	60965	2.48	ng/ul	0.00
9) 2-Chlorophenol-d4	7.49	132	25777	0.94	ng/ul	0.00
13) 4-Methylphenol-d8	8.66	113	19096	0.67	ng/ul	0.00
19) Nitrobenzene-d5	9.13	128	30899	2.37	ng/ul	0.00
22) 2-Nitrophenol-d4	9.85	143	4434	0.33	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.39	165	13025	0.48	ng/ul	0.00
29) 4-Chloroaniline-d4	10.90	131	238	0.01	ng/ul	0.00
44) Dimethylphthalate-d6	13.99	166	51865	0.65	ng/ul	0.00
47) Acenaphthylene-d8	14.29	160	225159	2.34	ng/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
58) Fluorene-d10	15.58	176	163970	2.38	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
71) Anthracene-d10	17.43	188	255435	2.61	ng/ul	0.00
79) Pyrene-d10	19.72	212	252070	3.68	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.72	264	152373	2.48	ng/ul	0.00

Target Compounds

					Ovalue	
70) Phenanthrene	17.37	178	302641	2.732	ng/ul	98
77) Fluoranthene	19.39	202	522871	4.589	ng/ul#	93
80) Pyrene	19.74	202	445012	5.213	ng/ul#	88
83) Benzo(a)anthracene	21.49	228	193814	2.406	ng/ul	99
85) Chrysene	21.54	228	187294	2.565	ng/ul	96
88) Benzo(b)fluoranthene	23.15	252	221406	2.859	ng/ul#	91
91) Benzo(a)pyrene	23.77	252	141957	1.942	ng/ul#	95
92) Indeno(1,2,3-cd)pyrene	26.30	276	97980	1.218	ng/ul#	85
94) Benzo(a,h,i)perylene	27.05	276	95892	1.472	ng/ul#	89

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

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