

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110817\
 Data File : BM012843.D
 Acq On : 09 Nov 2017 08:31
 Operator : SJ/JU
 Sample : I6096-02
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 B-76(1-3)-102317

Manual Integrations
 APPROVED

Sohil
 11/9/2017 3:49:06 PM

Quant Time: Nov 09 10:20:18 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM110417MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 09 05:48:58 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	151668	20.00	ng/ul	0.00
18) Naphthalene-d8	10.59	136	637442	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.44	164	390727	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.19	188	853465	20.00	ng/ul	0.00
77) Chrysene-d12	21.37	240	690061	20.00	ng/ul	0.00
85) Perylene-d12	23.66	264	760780	20.00	ng/ul	0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.29	96	5116	1.82	ng/uL	0.00
5) Phenol-d5	6.98	99	121355	10.75	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.13	67	60686	9.92	ng/ul	0.00
9) 2-Chlorophenol-d4	7.34	132	98445	10.47	ng/ul	0.00
13) 4-Methylphenol-d8	8.52	113	96995	9.96	ng/ul	0.00
19) Nitrobenzene-d5	8.96	128	46825	9.94	ng/ul	0.00
22) 2-Nitrophenol-d4	9.68	143	59670	11.07	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.22	165	124576	11.28	ng/ul	0.00
29) 4-Chloroaniline-d4	10.73	131	118786	11.14	ng/ul	0.00
43) Dimethylphthalate-d6	13.85	166	452372	13.99	ng/ul	0.00
46) Acenaphthylene-d8	14.13	160	436139	10.91	ng/ul	0.00
51) 4-Nitrophenol-d4	14.67	143	89828	17.34	ng/ul	0.00
57) Fluorene-d10	15.44	176	364471	13.16	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.58	200	13968	2.48	ng/ul	0.00
70) Anthracene-d10	17.29	188	778652	18.88	ng/ul	0.00
78) Pyrene-d10	19.59	212	904568	28.25	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.51	264	933186	25.94	ng/ul	0.01

Target Compounds

					Qvalue	
6) Phenol	7.01	94	12610	1.089	ng/ul#	91
44) Dimethylphthalate	13.90	163	180153	5.615	ng/ul	98
69) Phenanthrene	17.23	178	305205	6.543	ng/ul	99
71) Anthracene	17.32	178	67055	1.386	ng/ul	99
76) Fluoranthene	19.25	202	666195m	11.884	ng/ul	
79) Pyrene	19.62	202	659194	16.011	ng/ul#	92
82) Benzo(a)anthracene	21.36	228	362481	8.681	ng/ul	96
84) Chrysene	21.41	228	368682	9.742	ng/ul	97
87) Benzo(b)fluoranthene	22.97	252	589597	12.682	ng/ul#	95
88) Benzo(k)fluoranthene	23.01	252	192477m	4.355	ng/ul	
90) Benzo(a)pyrene	23.57	252	441369m	9.977	ng/ul	
91) Indeno(1,2,3-cd)pyrene	25.98	276	354936	6.657	ng/ul#	86
92) Dibenzo(a,h)anthracene	25.98	278	92982	2.060	ng/ul#	88
93) Benzo(g,h,i)perylene	26.69	276	317854	7.235	ng/ul#	90

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

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