

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121318\
 Data File : BM018150.D
 Acq On : 14 Dec 2018 04:00
 Operator : JU/SJ
 Sample : J6236-11
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F9E21

Manual Integrations
 APPROVED

Sohil
 12/14/2018 2:17:35 PM

Quant Time: Dec 14 05:35:26 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121318MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 14 05:30:21 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.51	152	140703	20.00	ng/ul	0.00
18) Naphthalene-d8	10.27	136	637925	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.16	164	345541	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.93	188	796511	20.00	ng/ul	0.00
77) Chrysene-d12	21.14	240	942436	20.00	ng/ul	0.00
85) Perylene-d12	23.30	264	947505	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.10	96	7893	2.68	ng/uL	0.00
5) Phenol-d5	6.70	99	171395	12.68	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.86	67	111638	15.39	ng/ul	0.00
9) 2-Chlorophenol-d4	7.05	132	146239	13.68	ng/ul	0.00
13) 4-Methylphenol-d8	8.23	113	125479	11.44	ng/ul	0.00
19) Nitrobenzene-d5	8.67	128	72021	13.74	ng/ul	0.00
22) 2-Nitrophenol-d4	9.38	143	86195	13.90	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.92	165	139507m	13.01	ng/ul	0.00
29) 4-Chloroaniline-d4	10.44	131	164172	17.33	ng/ul	0.00
43) Dimethylphthalate-d6	13.59	166	471029	15.41	ng/ul	0.00
46) Acenaphthylene-d8	13.86	160	518216	13.91	ng/ul	0.00
51) 4-Nitrophenol-d4	14.42	143	59587m	10.80	ng/ul	0.00
57) Fluorene-d10	15.17	176	396624	15.62	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.32	200	73492	12.11	ng/ul	0.00
70) Anthracene-d10	17.02	188	581266	14.44	ng/ul	0.00
78) Pyrene-d10	19.33	212	664695	14.71	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.17	264	737172	14.05	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.73	94	32057	2.411 ng/ul	95
44) Dimethylphthalate	13.64	163	65408	2.247 ng/ul	98
47) Acenaphthylene	13.89	152	65160	1.848 ng/ul	98
76) Fluoranthene	19.00	202	579791	10.750 ng/ul	100
79) Pyrene	19.36	202	633655	11.238 ng/ul	99
82) Benzo(a)anthracene	21.13	228	434903	7.291 ng/ul	99
84) Chrysene	21.18	228	420739	7.560 ng/ul	97
87) Benzo(b)fluoranthene	22.66	252	751375	13.307 ng/ul	98
88) Benzo(k)fluoranthene	22.70	252	275493m	4.903 ng/ul	
90) Benzo(a)pyrene	23.21	252	354745	6.369 ng/ul	99
91) Indeno(1,2,3-cd)pyrene	25.43	276	259811	4.111 ng/ul	96
92) Dibenzo(a,h)anthracene	25.43	278	70745m	1.317 ng/ul	
93) Benzo(g,h,i)perylene	26.09	276	192082	3.647 ng/ul	98

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

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