

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101218\
 Data File : BN003108.D
 Acq On : 11 Oct 2018 18:16
 Operator : MJ/SJ
 Sample : J5368-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A3Z47

Quant Time: Oct 11 18:52:40 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN100918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 11 16:42:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1863	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	7831	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	4453	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.11	188	350728	0.40	ng/ul	0.08
16) Chrysene-d12	21.37	240	85	0.40	ng/ul	0.14
20) Perylene-d12	23.69	264	1820	0.40	ng/ul	0.23
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	2672	0.22	ng/ul	-0.02
14) Fluoranthene-d10	0.00	212	0	0.00	ng/ul	
Target Compounds				Qvalue		
3) Naphthalene	10.44	128	2169	0.108	ng/ul	95
5) 2-Methylnaphthalene	12.08	142	1278	0.092	ng/ul	99
7) Acenaphthylene	13.98	152	3291	0.162	ng/ul#	35
8) Acenaphthene	14.32	153	4432	0.287	ng/ul	98
9) Fluorene	15.32	166	5149	0.291	ng/ul	95
12) Phenanthrene	17.15	178	34673	0.035	ng/ul	97
17) Pyrene	19.45	202	189753	574.772	ng/ul	96
18) Benzo(a)anthracene	21.26	228	109494	400.299	ng/ul	96
19) Chrysene	21.38	228	9927	35.741	ng/ul#	59
21) Benzo(b)fluoranthene	22.94	252	26714	4.076	ng/ul	85
22) Benzo(k)fluoranthene	23.13	252	2540	0.374	ng/ul#	1
23) Benzo(a)pyrene	23.58	252	1052	0.179	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	25.98	276	6841	1.013	ng/ul#	38
25) Dibenzo(a,h)anthracene	25.98	278	24207	4.480	ng/ul	99
26) Benzo(g,h,i)perylene	26.68	276	18742	3.293	ng/ul#	91

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

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