

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110518\
 Data File : BN003431.D
 Acq On : 05 Nov 2018 16:00
 Operator : MJ/SJ
 Sample : J5635-13
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A40J6

Quant Time: Nov 05 17:04:25 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN110118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 05 15:40:40 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	10791	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	45578	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	26376	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.33	188	3128968	0.40	ng/ul	0.10
16) Chrysene-d12	21.54	240	709	0.40	ng/ul	0.13
20) Perylene-d12	23.94	264	12	0.40	ng/ul	0.22
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	22238	0.30	ng/ul	0.00
14) Fluoranthene-d10	0.00	212	0	0.00	ng/ul	
Target Compounds						
					Qvalue	
3) Naphthalene	10.70	128	2889	0.023	ng/ul#	89
5) 2-Methylnaphthalene	12.31	142	4602	0.050	ng/ul	99
7) Acenaphthylene	14.21	152	4472	0.039	ng/ul#	85
17) Pyrene	19.65	202	32490	12.598	ng/ul	98
18) Benzo(a)anthracene	21.56	228	1692	0.761	ng/ul#	10
19) Chrysene	21.59	228	2310	0.914	ng/ul#	55
21) Benzo(b)fluoranthene	23.25	252	519	10.037	ng/ul#	1
22) Benzo(k)fluoranthene	23.25	252	519	10.363	ng/ul#	1
23) Benzo(a)pyrene	23.87	252	150	3.474	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	26.27	276	34	0.736	ng/ul#	32
25) Dibenzo(a,h)anthracene	26.36	278	1543	40.955	ng/ul#	1
26) Benzo(g,h,i)perylene	27.04	276	67	1.690	ng/ul#	1

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

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