

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110518\
 Data File : BN003430.D
 Acq On : 05 Nov 2018 15:25
 Operator : MJ/SJ
 Sample : J5635-12
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A40J5

Quant Time: Nov 05 16:08:34 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	10929	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	45536	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	25868	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.33	188	2796167	0.40	ng/ul	0.10
16) Chrysene-d12	21.54	240	2036	0.40	ng/ul	0.13
20) Perylene-d12	23.87	264	9170	0.40	ng/ul	0.15
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	20375	0.28	ng/ul	0.00
14) Fluoranthene-d10	19.37	212	947	0.00	ng/ul	0.12
Target Compounds				Qvalue		
3) Naphthalene	10.70	128	29636	0.232	ng/ul	99
5) 2-Methylnaphthalene	12.31	142	13766	0.150	ng/ul	100
7) Acenaphthylene	14.21	152	148268	1.319	ng/ul	99
8) Acenaphthene	14.55	153	9152	0.091	ng/ul	97
9) Fluorene	15.54	166	34478	0.292	ng/ul	99
12) Phenanthrene	17.36	178	382074	0.042	ng/ul	99
17) Pyrene	19.65	202	3917580	528.975	ng/ul	95
18) Benzo(a)anthracene	21.45	228	1762898	276.172	ng/ul	92
19) Chrysene	21.56	228	238140	32.825	ng/ul#	61
21) Benzo(b)fluoranthene	23.20	252	410486	10.389	ng/ul	90
22) Benzo(k)fluoranthene	23.20	252	410486	10.726	ng/ul#	90
23) Benzo(a)pyrene	23.77	252	396905	12.029	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	26.08	276	1047499	29.656	ng/ul	93
25) Dibenzo(a,h)anthracene	26.36	278	250889	8.714	ng/ul	95
26) Benzo(g,h,i)perylene	26.81	276	669147	22.091	ng/ul	99

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

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