

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

Manual Integrations
 APPROVED

Sohil
 11/6/2018 1:23:53 PM

Quant Time: Nov 05 17:20:30 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	17.23	188	55588m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	38669m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	33718m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	22238	0.30	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	49111m	0.31	ng/ul	0.00
Target Compounds				Qvalue		
5) 2-Methylnaphthalene	12.31	142	4602	0.050	ng/ul	99
12) Phenanthrene	17.28	178	9220	0.051	ng/ul	96
15) Fluoranthene	19.29	202	36436m	0.178	ng/ul	
17) Pyrene	19.65	202	32490	0.231	ng/ul	98
18) Benzo(a)anthracene	21.39	228	13600m	0.112	ng/ul	
19) Chrysene	21.45	228	15119m	0.110	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	20278m	0.140	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	7225m	0.051	ng/ul	
23) Benzo(a)pyrene	23.63	252	13030m	0.107	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.08	276	9913m	0.076	ng/ul	
26) Benzo(g,h,i)perylene	26.80	276	6807m	0.061	ng/ul	

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

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