

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110518\
 Data File : BN003427.D
 Acq On : 05 Nov 2018 13:41
 Operator : MJ/SJ
 Sample : J5704-07
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A40H1

Quant Time: Nov 05 14:40:53 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 13:38:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	9860	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	40925	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	23711	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.33	188	2691848	0.40	ng/ul	0.10
16) Chrysene-d12	21.54	240	1254	0.40	ng/ul	0.14
20) Perylene-d12	23.99	264	265	0.40	ng/ul	0.27
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	19628	0.30	ng/ul	0.00
14) Fluoranthene-d10	0.00	212	0	0.00	ng/ul	
Target Compounds				Qvalue		
8) Acenaphthene	14.55	153	1959	0.021	ng/ul	98
9) Fluorene	15.54	166	2455	0.023	ng/ul#	95
17) Pyrene	19.65	202	87271	19.132	ng/ul	99
18) Benzo(a)anthracene	21.56	228	4900	1.246	ng/ul#	11
19) Chrysene	21.56	228	4900	1.097	ng/ul#	16
21) Benzo(b)fluoranthene	23.33	252	376	0.329	ng/ul#	1
22) Benzo(k)fluoranthene	23.33	252	362	0.327	ng/ul#	1
23) Benzo(a)pyrene	23.87	252	506	0.531	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	26.35	276	1056	1.035	ng/ul#	73
25) Dibenzo(a,h)anthracene	26.35	278	4251	5.109	ng/ul#	3
26) Benzo(g,h,i)perylene	27.02	276	316	0.361	ng/ul#	1

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

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