

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110518\
 Data File : BN003429.D
 Acq On : 05 Nov 2018 14:51
 Operator : MJ/SJ
 Sample : J5704-23
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :

Quant Time: Nov 05 15:30:30 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 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	10439	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	43002	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	24531	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.33	188	2627235	0.40	ng/ul	0.10
16) Chrysene-d12	21.54	240	74	0.40	ng/ul	0.13
20) Perylene-d12	23.90	264	19	0.40	ng/ul	0.18
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	20569	0.30	ng/ul	0.00
14) Fluoranthene-d10	0.00	212	0	0.00	ng/ul	
Target Compounds				Qvalue		
7) Acenaphthylene	14.21	152	2140	0.020	ng/ul#	81
8) Acenaphthene	14.55	153	1907	0.020	ng/ul	99
17) Pyrene	19.65	202	72746	270.254	ng/ul	96
18) Benzo(a)anthracene	21.57	228	5805	25.021	ng/ul#	11
19) Chrysene	21.57	228	5805	22.015	ng/ul#	16
21) Benzo(b)fluoranthene	23.20	252	8453	103.248	ng/ul#	12
22) Benzo(k)fluoranthene	23.20	252	8453	106.601	ng/ul#	12
23) Benzo(a)pyrene	23.77	252	14753	215.792	ng/ul#	43
24) Indeno(1,2,3-cd)pyrene	26.20	276	953	13.022	ng/ul#	27
25) Dibenzo(a,h)anthracene	26.36	278	6131	102.778	ng/ul#	16
26) Benzo(g,h,i)perylene	27.03	276	1118	17.814	ng/ul#	1

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

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