

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100918\
 Data File : BN003019.D
 Acq On : 09 Oct 2018 10:10
 Operator : MJ/SJ
 Sample : J5115-13
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 JLC46

Quant Time: Oct 09 10:52:48 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN100918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 09 08:56:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1081	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	4744	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	3015	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	400459	0.40	ng/ul	0.08
16) Chrysene-d12	21.39	240	73	0.40	ng/ul	0.15
20) Perylene-d12	23.55	264	400	0.40	ng/ul	0.10
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	2273	0.31	ng/ul	-0.02
14) Fluoranthene-d10	0.00	212	0	0.00	ng/ul	
Target Compounds						
					Qvalue	
3) Naphthalene	10.44	128	1026	0.084	ng/ul#	87
5) 2-Methylnaphthalene	12.08	142	839	0.100	ng/ul	97
7) Acenaphthylene	13.98	152	1700	0.123	ng/ul#	90
8) Acenaphthene	14.32	153	539	0.051	ng/ul#	93
9) Fluorene	15.32	166	593	0.049	ng/ul#	92
17) Pyrene	19.45	202	96864	341.637	ng/ul	96
18) Benzo(a)anthracene	21.26	228	129034	549.280	ng/ul	97
19) Chrysene	21.47	228	370	1.551	ng/ul#	1
21) Benzo(b)fluoranthene	22.78	252	177706	123.360	ng/ul	89
22) Benzo(k)fluoranthene	22.95	252	8471	5.672	ng/ul#	1
23) Benzo(a)pyrene	23.49	252	13353	10.343	ng/ul#	9
24) Indeno(1,2,3-cd)pyrene	25.90	276	1752	1.181	ng/ul#	22
25) Dibenzo(a,h)anthracene	25.76	278	2509	2.113	ng/ul#	1
26) Benzo(g,h,i)perylene	26.33	276	55935	44.712	ng/ul	97

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

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