

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101218\
 Data File : BN003111.D
 Acq On : 11 Oct 2018 20:01
 Operator : MJ/SJ
 Sample : J5368-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :

Quant Time: Oct 12 06:38:31 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 : Fri Oct 12 05:29:36 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1649	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	6857	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.26	164	3681	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	7121m	0.40	ng/ul	-0.01
16) Chrysene-d12	21.23	240	4639m	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	4397m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	2531	0.24	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	4719m	0.23	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	803	0.046	ng/ul#	81
8) Acenaphthene	14.32	153	627	0.049	ng/ul	94
9) Fluorene	15.32	166	896	0.061	ng/ul#	96
12) Phenanthrene	17.06	178	7202	0.363	ng/ul	99
13) Anthracene	17.15	178	2058m	0.110	ng/ul	
15) Fluoranthene	19.09	202	5400m	0.232	ng/ul	
17) Pyrene	19.45	202	4070m	0.226	ng/ul	
18) Benzo(a)anthracene	21.22	228	1597	0.107	ng/ul#	89
19) Chrysene	21.27	228	1609	0.106	ng/ul#	87
21) Benzo(b)fluoranthene	22.78	252	1834m	0.116	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	820m	0.050	ng/ul	
23) Benzo(a)pyrene	23.34	252	1190	0.084	ng/ul#	24
24) Indeno(1,2,3-cd)pyrene	25.69	276	814	0.050	ng/ul#	94
26) Benzo(g,h,i)perylene	26.36	276	773	0.056	ng/ul#	36

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

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