

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090420\
 Data File : BN011709.D
 Acq On : 04 Sep 2020 19:20
 Operator : CG/JU
 Sample : L3840-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :

Quant Time: Sep 05 01:51:28 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN081220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 05 01:37:48 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	2542	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	9981	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	5766	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.13	188	12143m	0.40	ng/ul	0.00
16) Chrysene-d12	21.32	240	10534	0.40	ng/ul	0.00
20) Perylene-d12	23.57	264	10472	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	3019	0.21	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	7198	0.21	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.57	128	1159	0.040	ng/ul#	89
7) Acenaphthylene	14.10	152	2800	0.096	ng/ul	98
12) Phenanthrene	17.16	178	1752	0.043	ng/ul	96
13) Anthracene	17.26	178	2432m	0.072	ng/ul	
15) Fluoranthene	19.19	202	6169	0.131	ng/ul	96
17) Pyrene	19.55	202	11767m	0.286	ng/ul	
18) Benzo(a)anthracene	21.30	228	8119	0.230	ng/ul	95
19) Chrysene	21.36	228	14901	0.313	ng/ul	99
21) Benzo(b)fluoranthene	22.90	252	32894m	0.723	ng/ul	
22) Benzo(k)fluoranthene	22.94	252	13453m	0.267	ng/ul	
23) Benzo(a)pyrene	23.47	252	20815	0.544	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.84	276	12958	0.251	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.84	278	3910	0.093	ng/ul#	84
26) Benzo(g,h,i)perylene	26.53	276	10997	0.227	ng/ul	96

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

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