

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090420\
 Data File : BN011715.D
 Acq On : 05 Sep 2020 00:11
 Operator : CG/JU
 Sample : L3840-05
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :

Quant Time: Sep 05 02:08:01 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 01:53:42 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	2789	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	10746	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	6152	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.13	188	13223m	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	12124	0.40	ng/ul	0.00
20) Perylene-d12	23.56	264	12368	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	3352	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	7452	0.20	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.57	128	1916	0.061	ng/ul#	93
7) Acenaphthylene	14.10	152	4562	0.147	ng/ul	98
12) Phenanthrene	17.16	178	2632	0.059	ng/ul	98
13) Anthracene	17.26	178	5294	0.144	ng/ul	100
15) Fluoranthene	19.19	202	13984	0.272	ng/ul	98
17) Pyrene	19.55	202	19891m	0.420	ng/ul	
18) Benzo(a)anthracene	21.30	228	14990	0.369	ng/ul	97
19) Chrysene	21.35	228	33179	0.605	ng/ul	99
21) Benzo(b)fluoranthene	22.89	252	44163m	0.822	ng/ul	
22) Benzo(k)fluoranthene	22.93	252	18778m	0.316	ng/ul	
23) Benzo(a)pyrene	23.47	252	28236	0.625	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.82	276	20292	0.332	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.83	278	6181	0.124	ng/ul#	90
26) Benzo(g,h,i)perylene	26.51	276	17514	0.306	ng/ul	98

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

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