

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090420\
 Data File : BN011708.D
 Acq On : 04 Sep 2020 18:44
 Operator : CG/JU
 Sample : L3840-12
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :

Quant Time: Sep 05 01:48:27 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	2088	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	8076	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	4805	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.13	188	10369m	0.40	ng/ul	0.00
16) Chrysene-d12	21.32	240	9530	0.40	ng/ul	0.00
20) Perylene-d12	23.57	264	10028	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	4640	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	9960	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.57	128	2566	0.109	ng/ul	96
5) 2-Methylnaphthalene	12.19	142	450	0.030	ng/ul	97
7) Acenaphthylene	14.10	152	3143	0.130	ng/ul	97
9) Fluorene	15.43	166	610	0.027	ng/ul#	96
12) Phenanthrene	17.16	178	2784	0.079	ng/ul	96
13) Anthracene	17.26	178	4634	0.161	ng/ul	99
15) Fluoranthene	19.19	202	5315	0.132	ng/ul	93
17) Pyrene	19.55	202	21248m	0.570	ng/ul	
18) Benzo(a)anthracene	21.30	228	9177	0.287	ng/ul	96
19) Chrysene	21.36	228	23050	0.535	ng/ul	100
21) Benzo(b)fluoranthene	22.90	252	30683m	0.704	ng/ul	
22) Benzo(k)fluoranthene	22.94	252	13868m	0.288	ng/ul	
23) Benzo(a)pyrene	23.47	252	21147	0.577	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.83	276	15150	0.306	ng/ul	98
25) Dibenzo(a,h)anthracene	25.85	278	4463	0.111	ng/ul#	87
26) Benzo(a,h,i)perylene	26.53	276	14662	0.316	ng/ul	98

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

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