

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090820\
 Data File : BN011744.D
 Acq On : 08 Sep 2020 19:30
 Operator : CG/JU
 Sample : L3840-07
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AE0

Manual Integrations
 APPROVED

Sohil
 9/9/2020 3:46:17 PM

Quant Time: Sep 09 06:47:02 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN090820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 09 06:29:50 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	2044	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	7787	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	4435	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	9277m	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	7936	0.40	ng/ul	0.00
20) Perylene-d12	23.56	264	8101	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	3027	0.25	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	6592	0.27	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.57	128	1931	0.084	ng/ul	95
5) 2-Methylnaphthalene	12.19	142	331	0.022	ng/ul	96
7) Acenaphthylene	14.09	152	2327	0.109	ng/ul	95
12) Phenanthrene	17.16	178	1646	0.053	ng/ul#	91
13) Anthracene	17.25	178	2565m	0.092	ng/ul	
15) Fluoranthene	19.18	202	4439	0.128	ng/ul	95
17) Pyrene	19.54	202	10569	0.300	ng/ul#	94
18) Benzo(a)anthracene	21.29	228	6168	0.200	ng/ul	94
19) Chrysene	21.35	228	14464	0.420	ng/ul	99
21) Benzo(b)fluoranthene	22.88	252	22330m	0.588	ng/ul	
22) Benzo(k)fluoranthene	22.93	252	9534m	0.245	ng/ul	
23) Benzo(a)pyrene	23.46	252	14800	0.452	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.82	276	10128	0.235	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.82	278	3150	0.090	ng/ul#	82
26) Benzo(g,h,i)perylene	26.51	276	8751	0.232	ng/ul	94

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

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