

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090820\
 Data File : BN011747.D
 Acq On : 08 Sep 2020 21:19
 Operator : CG/JU
 Sample : L3840-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AA3

Manual Integrations
 APPROVED

Sohil
 9/9/2020 3:46:27 PM

Quant Time: Sep 09 06:56:48 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	7.74	152	1930	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	7608	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	4410	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	9231m	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	7808	0.40	ng/ul	0.00
20) Perylene-d12	23.56	264	7943	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	3631	0.30	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	7363	0.30	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.57	128	1397	0.062	ng/ul#	88
7) Acenaphthylene	14.09	152	1764	0.083	ng/ul#	93
12) Phenanthrene	17.16	178	1705	0.055	ng/ul#	91
13) Anthracene	17.25	178	2035	0.074	ng/ul#	92
15) Fluoranthene	19.18	202	4230	0.123	ng/ul	95
17) Pyrene	19.54	202	9507	0.274	ng/ul#	95
18) Benzo(a)anthracene	21.29	228	5578	0.184	ng/ul	94
19) Chrysene	21.34	228	11398	0.336	ng/ul	98
21) Benzo(b)fluoranthene	22.88	252	18467m	0.496	ng/ul	
22) Benzo(k)fluoranthene	22.93	252	7424m	0.194	ng/ul	
23) Benzo(a)pyrene	23.46	252	11215	0.349	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.82	276	7413	0.176	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.82	278	2164	0.063	ng/ul#	69
26) Benzo(g,h,i)perylene	26.51	276	6669	0.181	ng/ul#	92

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

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