

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090820\
 Data File : BN011754.D
 Acq On : 09 Sep 2020 02:08
 Operator : CG/JU
 Sample : L3843-05
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AD0

Manual Integrations
 APPROVED

Sohil
 9/9/2020 3:46:35 PM

Quant Time: Sep 09 07:35:17 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 07:05:35 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	2049	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	8074	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	4784	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	10291m	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	8956	0.40	ng/ul	0.00
20) Perylene-d12	23.56	264	9135	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	3556	0.28	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	7712	0.29	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.57	128	1600	0.067	ng/ul#	93
7) Acenaphthylene	14.09	152	3445	0.149	ng/ul	96
8) Acenaphthene	14.44	153	434	0.023	ng/ul	93
9) Fluorene	15.43	166	566	0.026	ng/ul#	95
12) Phenanthrene	17.16	178	3372	0.097	ng/ul#	96
13) Anthracene	17.25	178	4516	0.147	ng/ul	95
15) Fluoranthene	19.18	202	23955	0.623	ng/ul	98
17) Pyrene	19.54	202	42000	1.056	ng/ul	98
18) Benzo(a)anthracene	21.29	228	11908	0.342	ng/ul	95
19) Chrysene	21.34	228	21317	0.548	ng/ul	99
21) Benzo(b)fluoranthene	22.88	252	38959m	0.910	ng/ul	
22) Benzo(k)fluoranthene	22.92	252	13332m	0.304	ng/ul	
23) Benzo(a)pyrene	23.46	252	20639	0.558	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.82	276	12533	0.258	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.82	278	3664	0.092	ng/ul#	81
26) Benzo(a,h,i)perylene	26.51	276	11115	0.262	ng/ul	95

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

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