

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090820\
 Data File : BN011753.D
 Acq On : 09 Sep 2020 01:32
 Operator : CG/JU
 Sample : L3843-01
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AC0

Manual Integrations
 APPROVED

Sohil
 9/9/2020 3:46:34 PM

Quant Time: Sep 09 07:33:31 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	2174	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	8489	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	5070	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	10595m	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	9325	0.40	ng/ul	0.00
20) Perylene-d12	23.55	264	9473	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	4524	0.34	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	9069	0.33	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.57	128	1540	0.061	ng/ul#	89
7) Acenaphthylene	14.09	152	1554	0.064	ng/ul#	92
9) Fluorene	15.43	166	584	0.026	ng/ul#	92
12) Phenanthrene	17.16	178	3624	0.102	ng/ul	97
13) Anthracene	17.25	178	4435m	0.140	ng/ul	
15) Fluoranthene	19.18	202	31447	0.794	ng/ul	99
17) Pyrene	19.54	202	48475	1.171	ng/ul	98
18) Benzo(a)anthracene	21.29	228	20763	0.573	ng/ul	99
19) Chrysene	21.34	228	33248	0.821	ng/ul	98
21) Benzo(b)fluoranthene	22.88	252	31053m	0.699	ng/ul	
22) Benzo(k)fluoranthene	22.92	252	14240m	0.313	ng/ul	
23) Benzo(a)pyrene	23.46	252	20518	0.535	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.81	276	10293	0.204	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.82	278	3152	0.077	ng/ul#	78
26) Benzo(g,h,i)perylene	26.50	276	9876	0.224	ng/ul	96

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

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