

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121220\
 Data File : BN013098.D
 Acq On : 12 Dec 2020 19:10
 Operator : CG/JU
 Sample : L5063-05DL 5X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AD7DL

Manual Integrations
 APPROVED

mohammad
 12/15/2020 12:17:14 PM

Quant Time: Dec 14 01:35:02 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN120320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 14 00:25:18 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.37	152	843	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.13	136	3537	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.03	164	2321	0.40	ng/ul	-0.01
10) Phenanthrene-d10	16.79	188	4324	0.40	ng/ul	-0.02
16) Chrysene-d12	21.02	240	3299	0.40	ng/ul	-0.01
20) Perylene-d12	23.11	264	3484	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	215	0.04	ng/ul	-0.02
14) Fluoranthene-d10	18.84	212	440	0.04	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.17	128	8856	0.865	ng/ul	96
5) 2-Methylnaphthalene	11.80	142	40022	5.953	ng/ul	99
7) Acenaphthylene	13.75	152	2374	0.216	ng/ul#	38
8) Acenaphthene	14.09	153	2298	0.255	ng/ul	94
9) Fluorene	15.09	166	3949	0.382	ng/ul#	89
12) Phenanthrene	16.83	178	54359	3.749	ng/ul	98
13) Anthracene	16.92	178	3833m	0.311	ng/ul	
15) Fluoranthene	18.87	202	20461	1.188	ng/ul	98
17) Pyrene	19.23	202	29179	1.873	ng/ul	99
18) Benzo(a)anthracene	21.00	228	9562	0.701	ng/ul#	75
19) Chrysene	21.05	228	10257m	0.680	ng/ul	
21) Benzo(b)fluoranthene	22.50	252	13101m	0.813	ng/ul	
22) Benzo(k)fluoranthene	22.54	252	3944m	0.221	ng/ul	
23) Benzo(a)pyrene	23.02	252	8626	0.577	ng/ul#	75
24) Indeno(1,2,3-cd)pyrene	25.16	276	5952	0.303	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.16	278	1800	0.116	ng/ul#	1
26) Benzo(a,h,i)perylene	25.79	276	5833	0.328	ng/ul#	75

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

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