

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120320\
 Data File : BN012861.D
 Acq On : 03 Dec 2020 23:44
 Operator : CG/JU
 Sample : L4865-05DL 5X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0B77DL

Manual Integrations
 APPROVED

mohammad
 12/7/2020 8:51:44 AM

Quant Time: Dec 04 00:58:36 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 : Fri Dec 04 00:14:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	645	0.40	ng/ul	0.00
2) Naphthalene-d8	10.17	136	2481	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.06	164	1426	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.82	188	3056	0.40	ng/ul	0.00
16) Chrysene-d12	21.04	240	2704	0.40	ng/ul	0.00
20) Perylene-d12	23.14	264	2785	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.78	152	115	0.03	ng/ul	0.00
14) Fluoranthene-d10	18.86	212	264	0.03	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.21	128	159	0.022	ng/ul#	44
7) Acenaphthylene	13.78	152	152	0.023	ng/ul#	42
8) Acenaphthene	14.12	153	646	0.117	ng/ul	98
9) Fluorene	15.12	166	740	0.116	ng/ul#	98
12) Phenanthrene	16.86	178	27991	2.732	ng/ul	97
13) Anthracene	16.95	178	2719	0.312	ng/ul	98
15) Fluoranthene	18.89	202	41295	3.392	ng/ul	94
17) Pyrene	19.26	202	33747	2.642	ng/ul	96
18) Benzo(a)anthracene	21.02	228	12295	1.099	ng/ul	99
19) Chrysene	21.07	228	15766	1.275	ng/ul	98
21) Benzo(b)fluoranthene	22.52	252	18241m	1.416	ng/ul	
22) Benzo(k)fluoranthene	22.56	252	6691m	0.470	ng/ul	
23) Benzo(a)pyrene	23.05	252	10910	0.913	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.20	276	7922	0.504	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.20	278	1941	0.156	ng/ul#	87
26) Benzo(a,h,i)perylene	25.83	276	7129	0.502	ng/ul	98

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

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