

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120320\
 Data File : BN012857.D
 Acq On : 03 Dec 2020 21:20
 Operator : CG/JU
 Sample : L4866-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BC5

Manual Integrations
 APPROVED

mohammad
 12/7/2020 8:51:22 AM

Quant Time: Dec 04 00:37:46 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	571	0.40	ng/ul	0.00
2) Naphthalene-d8	10.16	136	2169	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.07	164	1313	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.82	188	2766m	0.40	ng/ul	0.00
16) Chrysene-d12	21.04	240	2278	0.40	ng/ul	0.00
20) Perylene-d12	23.14	264	2563	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.78	152	811	0.25	ng/ul	0.00
14) Fluoranthene-d10	18.86	212	1753	0.23	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.21	128	128	0.020	ng/ul#	34
5) 2-Methylnaphthalene	11.86	142	83	0.020	ng/ul	86
7) Acenaphthylene	13.78	152	166	0.027	ng/ul#	44
12) Phenanthrene	16.86	178	1190	0.128	ng/ul	93
13) Anthracene	16.95	178	215	0.027	ng/ul#	55
15) Fluoranthene	18.89	202	3413	0.310	ng/ul	100
17) Pyrene	19.26	202	3906	0.363	ng/ul#	92
18) Benzo(a)anthracene	21.03	228	1554	0.165	ng/ul	97
19) Chrysene	21.08	228	2009	0.193	ng/ul	97
21) Benzo(b)fluoranthene	22.52	252	2828m	0.239	ng/ul	
22) Benzo(k)fluoranthene	22.56	252	1138m	0.087	ng/ul	
23) Benzo(a)pyrene	23.05	252	2006	0.182	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	25.21	276	1535	0.106	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.21	278	345	0.030	ng/ul#	3
26) Benzo(g,h,i)perylene	25.84	276	1678	0.128	ng/ul#	85

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

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