

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120620\
 Data File : BN012927.D
 Acq On : 06 Dec 2020 20:53
 Operator : CG/JU
 Sample : L4859-07
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BD0

Manual Integrations
 APPROVED

mohammad
 12/7/2020 2:12:03 PM

Quant Time: Dec 07 04:02:39 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 07 03:06:53 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1827	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	7658	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.06	164	4357	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	9150	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	7455	0.40	ng/ul	0.00
20) Perylene-d12	23.14	264	6835	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	1441	0.13	ng/ul	-0.02
14) Fluoranthene-d10	18.86	212	3316	0.13	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.77	152	2645	0.128	ng/ul	100
8) Acenaphthene	14.12	153	815	0.048	ng/ul	98
9) Fluorene	15.12	166	1324	0.068	ng/ul	98
12) Phenanthrene	16.85	178	39301	1.281	ng/ul	97
13) Anthracene	16.95	178	5737	0.220	ng/ul	97
15) Fluoranthene	18.88	202	99249	2.722	ng/ul	99
17) Pyrene	19.25	202	81275	2.308	ng/ul	97
18) Benzo(a)anthracene	21.02	228	38663	1.254	ng/ul	99
19) Chrysene	21.07	228	39972	1.173	ng/ul	99
21) Benzo(b)fluoranthene	22.52	252	53932m	1.706	ng/ul	
22) Benzo(k)fluoranthene	22.55	252	19686m	0.563	ng/ul	
23) Benzo(a)pyrene	23.05	252	36231m	1.235	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.19	276	22732	0.590	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.20	278	5313	0.174	ng/ul	98
26) Benzo(g,h,i)perylene	25.82	276	19386	0.556	ng/ul	98

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

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