

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120720\
 Data File : BN012980.D
 Acq On : 08 Dec 2020 17:52
 Operator : CG/JU
 Sample : L4862-05DL 5X
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0B56DL

Manual Integrations
 APPROVED

mohammad
 12/9/2020 12:25:08 PM

Quant Time: Dec 09 00:29:15 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 : Wed Dec 09 00:11:42 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	1227	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	4902	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	2737	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	5454	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	3753	0.40	ng/ul	0.00
20) Perylene-d12	23.13	264	3368	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	294	0.04	ng/ul	0.00
14) Fluoranthene-d10	18.85	212	603	0.04	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.76	152	370	0.029	ng/ul#	73
12) Phenanthrene	16.85	178	5311	0.290	ng/ul	100
13) Anthracene	16.94	178	1312	0.084	ng/ul#	94
15) Fluoranthene	18.88	202	15735	0.724	ng/ul	96
17) Pyrene	19.25	202	12714	0.717	ng/ul	98
18) Benzo(a)anthracene	21.02	228	5263	0.339	ng/ul	99
19) Chrysene	21.07	228	5590	0.326	ng/ul	98
21) Benzo(b)fluoranthene	22.52	252	6527m	0.419	ng/ul	
22) Benzo(k)fluoranthene	22.55	252	2639m	0.153	ng/ul	
23) Benzo(a)pyrene	23.04	252	4103	0.284	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.19	276	2742	0.144	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.19	278	666	0.044	ng/ul#	42
26) Benzo(g,h,i)perylene	25.82	276	2373	0.138	ng/ul#	92

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

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