

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120720\
 Data File : BN012984.D
 Acq On : 08 Dec 2020 20:17
 Operator : CG/JU
 Sample : L4862-01DL 5X
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AY8DL

Manual Integrations
 APPROVED

mohammad
 12/9/2020 12:25:19 PM

Quant Time: Dec 09 00:45:24 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	1296	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	4985	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	2740	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.80	188	5524	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	3665	0.40	ng/ul	0.00
20) Perylene-d12	23.13	264	3366	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	399	0.05	ng/ul	0.00
14) Fluoranthene-d10	18.85	212	729	0.05	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.20	128	446m	0.031	ng/ul	
8) Acenaphthene	14.11	153	1166	0.110	ng/ul	98
9) Fluorene	15.11	166	603	0.049	ng/ul#	93
12) Phenanthrene	16.85	178	12003	0.648	ng/ul	97
13) Anthracene	16.94	178	2023	0.128	ng/ul	96
15) Fluoranthene	18.88	202	19371	0.880	ng/ul	95
17) Pyrene	19.25	202	17546	1.014	ng/ul	99
18) Benzo(a)anthracene	21.02	228	7121	0.470	ng/ul	98
19) Chrysene	21.07	228	8427	0.503	ng/ul	98
21) Benzo(b)fluoranthene	22.51	252	11270m	0.724	ng/ul	
22) Benzo(k)fluoranthene	22.55	252	3813m	0.222	ng/ul	
23) Benzo(a)pyrene	23.04	252	7368	0.510	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.19	276	5367	0.283	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.19	278	1203	0.080	ng/ul#	71
26) Benzo(g,h,i)perylene	25.81	276	5726	0.334	ng/ul	98

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

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