

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120720\
 Data File : BN013011.D
 Acq On : 09 Dec 2020 14:18
 Operator : CG/JU
 Sample : L4862-08DL 10X
 Misc :
 ALS Vial : 81 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0B62DL

Manual Integrations
 APPROVED

mohammad
 12/9/2020 5:39:35 PM

Quant Time: Dec 09 14:50:16 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 07:33:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	984	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.14	136	3984	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.05	164	2184	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	4519	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	3614	0.40	ng/ul	0.00
20) Perylene-d12	23.14	264	3441	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.75	152	226	0.04	ng/ul	0.00
14) Fluoranthene-d10	18.85	212	473	0.04	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.76	152	748	0.072	ng/ul#	89
9) Fluorene	15.11	166	314	0.032	ng/ul#	91
12) Phenanthrene	16.84	178	5706	0.377	ng/ul	99
13) Anthracene	16.94	178	1156	0.090	ng/ul#	92
15) Fluoranthene	18.88	202	15589	0.866	ng/ul	99
17) Pyrene	19.25	202	13246	0.776	ng/ul	99
18) Benzo(a)anthracene	21.02	228	6251	0.418	ng/ul	99
19) Chrysene	21.07	228	8057	0.488	ng/ul	98
21) Benzo(b)fluoranthene	22.52	252	10153m	0.638	ng/ul	
22) Benzo(k)fluoranthene	22.56	252	4154m	0.236	ng/ul	
23) Benzo(a)pyrene	23.04	252	6159	0.417	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.19	276	4286	0.221	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.20	278	1075	0.070	ng/ul#	58
26) Benzo(g,h,i)perylene	25.82	276	4086	0.233	ng/ul	97

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

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