

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120720\
 Data File : BN012946.D
 Acq On : 07 Dec 2020 19:19
 Operator : CG/JU
 Sample : L4859-03DL 2X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0B93DL

Manual Integrations
 APPROVED

mohammad
 12/9/2020 12:21:26 PM

Quant Time: Dec 08 00:35:25 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 : Tue Dec 08 00:15:39 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	877	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	3573	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	2024	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	4245	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	3596	0.40	ng/ul	0.00
20) Perylene-d12	23.14	264	3640	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	425	0.08	ng/ul	0.00
14) Fluoranthene-d10	18.86	212	986	0.09	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.20	128	299	0.029	ng/ul#	65
5) 2-Methylnaphthalene	11.84	142	169	0.025	ng/ul	96
7) Acenaphthylene	13.77	152	433	0.045	ng/ul#	79
8) Acenaphthene	14.12	153	179	0.023	ng/ul#	91
9) Fluorene	15.11	166	215	0.024	ng/ul#	92
12) Phenanthrene	16.85	178	6376	0.448	ng/ul	99
13) Anthracene	16.94	178	1053	0.087	ng/ul#	91
15) Fluoranthene	18.88	202	18041	1.067	ng/ul	98
17) Pyrene	19.25	202	17146	1.009	ng/ul	99
18) Benzo(a)anthracene	21.02	228	7661	0.515	ng/ul	99
19) Chrysene	21.07	228	9687	0.589	ng/ul	99
21) Benzo(b)fluoranthene	22.52	252	13266m	0.788	ng/ul	
22) Benzo(k)fluoranthene	22.55	252	4955m	0.266	ng/ul	
23) Benzo(a)pyrene	23.04	252	8468	0.542	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.19	276	5711	0.278	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.19	278	1348	0.083	ng/ul#	75
26) Benzo(a,h,i)perylene	25.82	276	5191	0.280	ng/ul	99

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

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