

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120720\
 Data File : BN012942.D
 Acq On : 07 Dec 2020 16:54
 Operator : CG/JU
 Sample : L4859-02DL 10X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0B26DL

Manual Integrations
 APPROVED

mohammad
 12/9/2020 12:21:14 PM

Quant Time: Dec 07 17:27:36 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 12:24:06 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	914	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	3701	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.06	164	2096	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	4416	0.40	ng/ul	0.00
16) Chrysene-d12	21.04	240	3751	0.40	ng/ul	0.00
20) Perylene-d12	23.14	264	3753	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.78	152	72	0.01	ng/ul	0.02
14) Fluoranthene-d10	18.86	212	147	0.01	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.77	152	290	0.029	ng/ul#	65
8) Acenaphthene	14.12	153	258	0.032	ng/ul	94
9) Fluorene	15.12	166	255	0.027	ng/ul#	88
12) Phenanthrene	16.85	178	6696	0.452	ng/ul	99
13) Anthracene	16.95	178	1115	0.088	ng/ul#	94
15) Fluoranthene	18.88	202	15659	0.890	ng/ul	97
17) Pyrene	19.25	202	14154	0.799	ng/ul	98
18) Benzo(a)anthracene	21.02	228	6713	0.433	ng/ul	98
19) Chrysene	21.07	228	7786	0.454	ng/ul	99
21) Benzo(b)fluoranthene	22.52	252	11091m	0.639	ng/ul	
22) Benzo(k)fluoranthene	22.56	252	4048m	0.211	ng/ul	
23) Benzo(a)pyrene	23.05	252	6993	0.434	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.20	276	4839	0.229	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.20	278	1152	0.069	ng/ul#	72
26) Benzo(g,h,i)perylene	25.82	276	4316	0.225	ng/ul	98

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

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