

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120720\
 Data File : BN012952.D
 Acq On : 07 Dec 2020 22:55
 Operator : CG/JU
 Sample : L4859-07DL 2X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BD0DL

Manual Integrations
 APPROVED

mohammad
 12/9/2020 12:21:39 PM

Quant Time: Dec 08 00:58:29 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	975	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	4074	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	2304	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	4720	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	3896	0.40	ng/ul	0.00
20) Perylene-d12	23.13	264	3901	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	458	0.08	ng/ul	0.00
14) Fluoranthene-d10	18.85	212	1016	0.08	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.76	152	950	0.087	ng/ul#	92
8) Acenaphthene	14.12	153	295	0.033	ng/ul	95
9) Fluorene	15.12	166	444	0.043	ng/ul#	96
12) Phenanthrene	16.85	178	12216	0.772	ng/ul	97
13) Anthracene	16.95	178	1818	0.135	ng/ul	97
15) Fluoranthene	18.88	202	31191	1.659	ng/ul	98
17) Pyrene	19.25	202	26979	1.466	ng/ul	100
18) Benzo(a)anthracene	21.02	228	12032	0.746	ng/ul	99
19) Chrysene	21.07	228	13192	0.741	ng/ul	99
21) Benzo(b)fluoranthene	22.52	252	18120m	1.004	ng/ul	
22) Benzo(k)fluoranthene	22.55	252	6492m	0.325	ng/ul	
23) Benzo(a)pyrene	23.04	252	12432m	0.742	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.19	276	7734	0.352	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.20	278	1775	0.102	ng/ul#	80
26) Benzo(g,h,i)perylene	25.82	276	6702	0.337	ng/ul	98

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

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