

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090420\
 Data File : BN011721.D
 Acq On : 05 Sep 2020 03:48
 Operator : CG/JU
 Sample : L3843-01
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :

Quant Time: Sep 05 06:28:15 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN081220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 05 05:40:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	5	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	95	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	524	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	600m	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	682	0.40	ng/ul	0.00
20) Perylene-d12	23.55	264	569m	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	5361	39.88	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	10933	6.48	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.57	128	1812	6.543	ng/ul#	94
5) 2-Methylnaphthalene	12.19	142	291	1.662	ng/ul	95
7) Acenaphthylene	14.10	152	1767	0.670	ng/ul	93
8) Acenaphthene	14.44	153	440	0.214	ng/ul	96
9) Fluorene	15.43	166	686	0.283	ng/ul#	93
12) Phenanthrene	17.16	178	4284	2.110	ng/ul	100
13) Anthracene	17.26	178	5200m	3.114	ng/ul	
15) Fluoranthene	19.18	202	37531	16.074	ng/ul	97
17) Pyrene	19.55	202	56844	21.324	ng/ul	98
18) Benzo(a)anthracene	21.30	228	24834	10.853	ng/ul	98
19) Chrysene	21.35	228	39815	12.902	ng/ul	99
21) Benzo(b)fluoranthene	22.89	252	37270m	15.070	ng/ul	
22) Benzo(k)fluoranthene	22.93	252	18023m	6.591	ng/ul	
23) Benzo(a)pyrene	23.47	252	24772	11.921	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.83	276	12561	4.470	ng/ul	99
25) Dibenzo(a,h)anthracene	25.83	278	3756	1.640	ng/ul#	84
26) Benzo(a,h,i)perylene	26.51	276	12024	4.572	ng/ul	97

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

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