

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101420\
 Data File : BN012218.D
 Acq On : 15 Oct 2020 05:57
 Operator : CG/JU
 Sample : L4184-15DL 5X
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :

Quant Time: Oct 15 07:59:02 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 15 04:47:44 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	2090	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	8758	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	4978	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	10701	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	8628	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	8886	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.99	152	516	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	1143	0.04	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.44	128	4817	0.191	ng/ul	99
5) 2-Methylnaphthalene	12.07	142	1409	0.088	ng/ul	98
7) Acenaphthylene	13.98	152	4371	0.202	ng/ul	98
8) Acenaphthene	14.32	153	2450	0.134	ng/ul	98
9) Fluorene	15.31	166	3327	0.164	ng/ul#	98
11) Pentachlorophenol	16.67	266	147	0.048	ng/ul	95
12) Phenanthrene	17.05	178	61187	1.794	ng/ul	98
13) Anthracene	17.14	178	14645	0.501	ng/ul	99
15) Fluoranthene	19.07	202	141855	3.556	ng/ul	99
17) Pyrene	19.44	202	122005	3.136	ng/ul	98
18) Benzo(a)anthracene	21.19	228	70050	2.192	ng/ul	98
19) Chrysene	21.24	228	67276	1.728	ng/ul	98
21) Benzo(b)fluoranthene	22.74	252	95588m	2.663	ng/ul	
22) Benzo(k)fluoranthene	22.78	252	32796m	0.826	ng/ul	
23) Benzo(a)pyrene	23.30	252	64105	1.836	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	25.58	276	48130	1.073	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.58	278	12705	0.354	ng/ul#	83
26) Benzo(g,h,i)perylene	26.25	276	45439	1.105	ng/ul	98

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

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