

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101420\
 Data File : BN012219.D
 Acq On : 15 Oct 2020 06:34
 Operator : CG/JU
 Sample : L4184-15
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :

Quant Time: Oct 15 08:03:21 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	7.62	152	1848	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	7532	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	4551	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	9354m	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	6670	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	6869	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.99	152	2696	0.24	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	5488	0.21	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	26190	1.205	ng/ul	97
5) 2-Methylnaphthalene	12.06	142	8079	0.585	ng/ul	99
7) Acenaphthylene	13.98	152	22408	1.135	ng/ul	99
8) Acenaphthene	14.32	153	12977	0.778	ng/ul	98
9) Fluorene	15.31	166	17655	0.950	ng/ul	98
11) Pentachlorophenol	16.66	266	786	0.291	ng/ul	99
12) Phenanthrene	17.05	178	302375	10.141	ng/ul	98
13) Anthracene	17.14	178	77492	3.030	ng/ul	98
15) Fluoranthene	19.08	202	657215	18.850	ng/ul	98
17) Pyrene	19.44	202	554773	18.448	ng/ul	99
18) Benzo(a)anthracene	21.20	228	333902	13.515	ng/ul	98
19) Chrysene	21.25	228	309900	10.296	ng/ul	97
21) Benzo(b)fluoranthene	22.75	252	455061m	16.402	ng/ul	
22) Benzo(k)fluoranthene	22.79	252	155538m	5.065	ng/ul	
23) Benzo(a)pyrene	23.31	252	311298	11.534	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	25.59	276	232938	6.718	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.60	278	62910	2.268	ng/ul#	91
26) Benzo(g,h,i)perylene	26.27	276	217634	6.848	ng/ul	98

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

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