

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN072120\
 Data File : BN011214.D
 Acq On : 21 Jul 2020 10:55
 Operator : CG/JU
 Sample : L3260-03
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 X2M73

Quant Time: Jul 21 11:49:21 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN072020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 21 11:43:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.49	152	1015	0.40	ng/ul	0.00
2) Naphthalene-d8	10.24	136	3527	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.13	164	2185	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.88	188	4829	0.40	ng/ul	0.00
16) Chrysene-d12	21.10	240	3666	0.40	ng/ul	0.00
20) Perylene-d12	23.23	264	3185	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.84	152	2037	0.24	ng/ul	0.00
14) Fluoranthene-d10	18.92	212	5542	0.27	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.29	128	3498	0.332	ng/ul	97
5) 2-Methylnaphthalene	11.92	142	4162	0.603	ng/ul	100
7) Acenaphthylene	13.84	152	2683	0.253	ng/ul	99
9) Fluorene	15.19	166	4906	0.519	ng/ul	98
11) Pentachlorophenol	16.54	266	1047	1.015	ng/ul	98
12) Phenanthrene	16.92	178	5263	0.331	ng/ul	99
13) Anthracene	17.01	178	6822	0.520	ng/ul	100
15) Fluoranthene	18.95	202	5549	0.312	ng/ul	99
17) Pyrene	19.32	202	7190	0.367	ng/ul	99
18) Benzo(a)anthracene	21.08	228	6322	0.488	ng/ul	98
19) Chrysene	21.13	228	8315	0.482	ng/ul	99
21) Benzo(b)fluoranthene	22.60	252	10721	0.779	ng/ul	91
22) Benzo(k)fluoranthene	22.65	252	5762	0.365	ng/ul	99
23) Benzo(a)pyrene	23.14	252	4249	0.367	ng/ul	100
25) Dibenzo(a,h)anthracene	25.33	278	8594	0.836	ng/ul#	90
26) Benzo(a,h,i)perylene	25.95	276	4490	0.331	ng/ul	98

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

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