

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM050120\
 Data File : BM025966.D
 Acq On : 01 May 2020 13:38
 Operator : MA/SJ
 Sample : L2321-03
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X2M11

Quant Time: May 01 14:10:08 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM041520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 01 12:24:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	1165	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	4040	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	2420	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	5718	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	4259	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	3943	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	2614	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	6519	0.40	ng/ul	0.00
Target Compounds						
3) Naphthalene	10.37	128	9200	0.72	ng/ul	98
5) 2-Methylnaphthalene	12.00	142	4337	0.52	ng/ul	98
9) Fluorene	15.26	166	7763	0.69	ng/ul	98
11) Pentachlorophenol	16.60	266	5969	4.13	ng/ul	96
13) Anthracene	17.08	178	15970	0.97	ng/ul	98
15) Fluoranthene	19.01	202	17630	0.78	ng/ul	100
17) Pyrene	19.37	202	19927	0.97	ng/ul	99
19) Chrysene	21.18	228	12838	0.64	ng/ul	99
21) Benzo(b)fluoranthene	22.65	252	10323	0.57	ng/ul	96
23) Benzo(a)pyrene	23.20	252	11544	0.77	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.42	278	20985	1.64	ng/ul#	90

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

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