

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM050120\
 Data File : BM025969.D
 Acq On : 01 May 2020 15:28
 Operator : MA/SJ
 Sample : L2321-08
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DFTPP02

Quant Time: May 01 16:06:30 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	1504	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	6075	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	3506	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	7958	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	6888	0.40	ng/ul	0.00
20) Perylene-d12	23.28	264	6716	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	2312	0.23	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	7654	0.34	ng/ul	0.00
Target Compounds						
3) Naphthalene	10.37	128	9127	0.48	ng/ul	99
5) 2-Methylnaphthalene	11.99	142	5978	0.47	ng/ul	98
9) Fluorene	15.25	166	9168	0.57	ng/ul	99
11) Pentachlorophenol	16.60	266	4491	2.23	ng/ul	95
12) Phenanthrene	16.98	178	17053	0.63	ng/ul	100
15) Fluoranthene	19.00	202	19663	0.62	ng/ul	97
18) Benzo(a)anthracene	21.12	228	21316	0.82	ng/ul	99
22) Benzo(k)fluoranthene	22.69	252	23924	0.71	ng/ul#	72
24) Indeno(1,2,3-cd)pyrene	25.40	276	15469	0.55	ng/ul#	94
26) Benzo(g,h,i)perylene	26.04	276	3148	0.12	ng/ul#	93

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

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