

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM050420\
 Data File : BM025971.D
 Acq On : 04 May 2020 14:02
 Operator : MA/SJ
 Sample : L2321-08
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X2M16

Quant Time: May 04 14:43:09 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 : Mon May 04 14:41:11 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	2227	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	8591	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	5107	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	11524	0.40	ng/ul	0.00
16) Chrysene-d12	21.13	240	9610	0.40	ng/ul	0.00
20) Perylene-d12	23.28	264	8499	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	3502	0.25	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	11060	0.33	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.37	128	13100	0.48	ng/ul	98
5) 2-Methylnaphthalene	11.99	142	8709	0.49	ng/ul	98
9) Fluorene	15.25	166	13288	0.56	ng/ul	98
11) Pentachlorophenol	16.60	266	6743	2.32	ng/ul	96
12) Phenanthrene	16.98	178	24731	0.63	ng/ul	98
15) Fluoranthene	19.00	202	26827	0.59	ng/ul	97
18) Benzo(a)anthracene	21.12	228	29262	0.80	ng/ul	98
22) Benzo(k)fluoranthene	22.69	252	27884	0.65	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	25.40	276	18506	0.52	ng/ul#	97
26) Benzo(g,h,i)perylene	26.03	276	18457	0.56	ng/ul	95

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

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