

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021821\
 Data File : BM028828.D
 Acq On : 19 Feb 2021 20:06
 Operator : CG/JU
 Sample : M1412-16
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBGY1

Quant Time: Feb 20 01:39:11 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM020521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Feb 20 01:10:53 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	341	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	1386	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	807	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	1615	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	1440	0.40	ng/ul	0.00
20) Perylene-d12	23.65	264	1439	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	390	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	855	0.21	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.74	128	138	0.040	ng/ul#	19
5) 2-Methylnaphthalene	12.35	142	53	0.023	ng/ul	94
9) Fluorene	15.58	166	60	0.021	ng/ul#	75
11) Pentachlorophenol	16.93	266	28	0.066	ng/ul#	29
12) Phenanthrene	17.31	178	388	0.086	ng/ul#	67
13) Anthracene	17.40	178	169	0.039	ng/ul#	33
15) Fluoranthene	19.32	202	423	0.082	ng/ul	70
17) Pyrene	19.68	202	324	0.062	ng/ul#	62
18) Benzo(a)anthracene	21.41	228	108	0.022	ng/ul#	1
19) Chrysene	21.46	228	184	0.038	ng/ul#	24
21) Benzo(b)fluoranthene	22.98	252	198	0.040	ng/ul#	1

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

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