

Data Path : Z:\HPCHEM1\BNA_M\Data\BM112216\
 Data File : BM007998.D
 Acq On : 22 Nov 2016 17:24
 Operator : UM/SJ
 Sample : H5622-23DL 5X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4WM0DL

Quant Time: Nov 22 17:56:42 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM112116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 22 16:05:09 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	599	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.50	136	1977	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.35	164	1436	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.09	188	2464	0.40	ng/ul	-0.02
16) Chrysene-d12	21.29	240	2625	0.40	ng/ul	-0.01
20) Perylene-d12	23.53	264	2451	0.40	ng/ul	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.09	152	224	0.08	ng/ul	-0.03
14) Fluoranthene-d10	19.13	212	608	0.09	ng/ul	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.55	128	4321	0.78	ng/ul	98
5) 2-Methylnaphthalene	12.16	142	1767	0.50	ng/ul	95
7) Acenaphthylene	14.06	152	156	0.02	ng/ul#	1
8) Acenaphthene	14.40	153	395	0.08	ng/ul#	81
9) Fluorene	15.41	166	879	0.15	ng/ul#	94
11) Pentachlorophenol	16.76	266	20	0.02	ng/ul#	63
12) Phenanthrene	17.14	178	1246	0.16	ng/ul#	84
13) Anthracene	17.23	178	206	0.03	ng/ul#	28
15) Fluoranthene	19.17	202	370	0.04	ng/ul#	1
17) Pyrene	19.52	202	338	0.04	ng/ul#	1
23) Benzo(a)pyrene	23.39	252	264	0.03	ng/ul#	1

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

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