

Data Path : Z:\HPCHEM1\BNA_M\Data\BM031017\
 Data File : BM009170.D
 Acq On : 10 Mar 2017 11:35
 Operator : SJ/MA
 Sample : I2107-16
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AL1

Quant Time: Mar 10 15:29:58 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM030317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 10 15:19:24 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	238	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	713	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	410	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.36	188	32498	0.40	ng/ul	0.10
16) Chrysene-d12	21.44	240	933	0.40	ng/ul	0.00
20) Perylene-d12	23.77	264	922	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.27	152	164	0.14	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	399	0.00	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.73	128	258	0.13	ng/ul#	56
5) 2-Methylnaphthalene	12.34	142	214	0.14	ng/ul	99
7) Acenaphthylene	14.23	152	50	0.02	ng/ul#	1
9) Fluorene	15.51	166	78	0.04	ng/ul#	50
17) Pyrene	19.69	202	791	0.25	ng/ul#	87
18) Benzo(a)anthracene	21.42	228	369	0.12	ng/ul#	75
19) Chrysene	21.47	228	461	0.16	ng/ul#	76
21) Benzo(b)fluoranthene	23.07	252	791	0.22	ng/ul#	30
22) Benzo(k)fluoranthene	23.07	252	779	0.24	ng/ul#	27
23) Benzo(a)pyrene	23.67	252	368	0.11	ng/ul#	57
24) Indeno(1,2,3-cd)pyrene	26.15	276	270	0.07	ng/ul	92
25) Dibenzo(a,h)anthracene	26.17	278	75	0.02	ng/ul#	1
26) Benzo(g,h,i)perylene	26.89	276	233	0.07	ng/ul#	43

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

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