

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM042417\
 Data File : BM009694.D
 Acq On : 24 Apr 2017 16:51
 Operator : SJ/MA
 Sample : I2779-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X2929

Quant Time: Apr 25 04:32:01 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM042417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 24 15:43:15 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	111	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	359	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	235	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	584	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	689	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	765	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.22	152	185	0.34	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	640	0.37	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.68	128	411	0.43	ng/ul#	66
5) 2-Methylnaphthalene	12.29	142	625	0.89	ng/ul	98
7) Acenaphthylene	14.19	152	573	0.50	ng/ul#	82
8) Acenaphthene	14.54	153	1139	1.25	ng/ul	95
11) Pentachlorophenol	16.88	266	255	1.13	ng/ul	93
12) Phenanthrene	17.26	178	1026	0.59	ng/ul#	94
15) Fluoranthene	19.28	202	2154	0.94	ng/ul	98
17) Pyrene	19.64	202	1807	0.79	ng/ul	96
18) Benzo(a)anthracene	21.39	228	1174	0.51	ng/ul	96
21) Benzo(b)fluoranthene	23.01	252	1998	0.76	ng/ul	91
22) Benzo(k)fluoranthene	23.06	252	2144	0.85	ng/ul#	92
23) Benzo(a)pyrene	23.61	252	1226	0.48	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	26.07	276	2208	0.76	ng/ul#	92
25) Dibenzo(a,h)anthracene	26.08	278	2219	0.92	ng/ul#	92
26) Benzo(g,h,i)perylene	26.79	276	2347	0.94	ng/ul#	91

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

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