

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE060817\
 Data File : BE093087.D
 Acq On : 8 Jun 2017 18:58
 Operator : SJ/MA
 Sample : MDL-S-4
 Misc : 0.07 PPM
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Jun 08 19:30:32 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 08 17:42:17 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	5167	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	22791	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	11431	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	31925	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	36405	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	33586	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	11701	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	29255	0.34	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.57	128	2920	0.052	ng/ul#	95
5) 2-Methylnaphthalene	12.19	142	1791	0.046	ng/ul	100
7) Acenaphthylene	14.09	152	2167	0.044	ng/ul	95
8) Acenaphthene	14.43	153	1931	0.049	ng/ul	98
9) Fluorene	15.42	166	2154	0.046	ng/ul	97
11) Pentachlorophenol	16.75	266	166	0.035	ng/ul	99
12) Phenanthrene	17.24	178	3201	0.036	ng/ul	96
13) Anthracene	17.24	178	3431	0.045	ng/ul	98
15) Fluoranthene	19.16	202	5065	0.047	ng/ul	85
17) Pyrene	19.52	202	5154	0.048	ng/ul#	85
18) Benzo(a)anthracene	21.24	228	4598	0.047	ng/ul#	85
19) Chrysene	21.29	228	5168	0.051	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	4369	0.043	ng/ul	91
22) Benzo(k)fluoranthene	22.98	252	5019	0.046	ng/ul#	89
23) Benzo(a)pyrene	23.56	252	4128	0.043	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	26.18	276	5042	0.045	ng/ul#	47
25) Dibenzo(a,h)anthracene	26.21	278	4157	0.044	ng/ul#	80
26) Benzo(g,h,i)perylene	26.96	276	4268	0.045	ng/ul	93

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

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