

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE060817\
 Data File : BE093084.D
 Acq On : 8 Jun 2017 17:07
 Operator : SJ/MA
 Sample : MDL-S-1
 Misc : 0.07 PPM
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Jun 08 17:42:03 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	7.74	152	4607	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	21027	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	10757	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	30058	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	34541	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	32404	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	10605	0.34	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	26864	0.33	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.57	128	2654	0.051	ng/ul#	94
5) 2-Methylnaphthalene	12.19	142	1663	0.047	ng/ul	98
7) Acenaphthylene	14.08	152	1918	0.042	ng/ul#	93
8) Acenaphthene	14.43	153	1827	0.049	ng/ul	95
9) Fluorene	15.42	166	1932	0.044	ng/ul	96
11) Pentachlorophenol	16.75	266	164	0.037	ng/ul	95
12) Phenanthrene	17.24	178	2980	0.036	ng/ul	97
13) Anthracene	17.24	178	3126	0.043	ng/ul	99
15) Fluoranthene	19.16	202	4657	0.046	ng/ul	84
17) Pyrene	19.52	202	4690	0.046	ng/ul#	86
18) Benzo(a)anthracene	21.25	228	4376	0.047	ng/ul#	85
19) Chrysene	21.30	228	4734	0.050	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	4013	0.041	ng/ul	93
22) Benzo(k)fluoranthene	22.98	252	4715	0.045	ng/ul#	87
23) Benzo(a)pyrene	23.56	252	4270	0.046	ng/ul	92
24) Indeno(1,2,3-cd)pyrene	26.19	276	4675	0.043	ng/ul#	85
25) Dibenzo(a,h)anthracene	26.21	278	4023	0.044	ng/ul#	81
26) Benzo(g,h,i)perylene	26.96	276	4014	0.043	ng/ul#	90

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

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