

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE060817\
 Data File : BE093086.D
 Acq On : 8 Jun 2017 18:22
 Operator : SJ/MA
 Sample : MDL-S-3
 Misc : 0.07 PPM
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Jun 08 18:53:02 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	4732	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	21462	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	10694	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	30028	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	33161	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	31539	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	10887	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	27054	0.33	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.57	128	2665	0.050	ng/ul#	96
5) 2-Methylnaphthalene	12.19	142	1704	0.047	ng/ul	97
7) Acenaphthylene	14.09	152	1957	0.043	ng/ul	96
8) Acenaphthene	14.43	153	1848	0.050	ng/ul	94
9) Fluorene	15.42	166	2019	0.047	ng/ul	96
11) Pentachlorophenol	16.75	266	179	0.040	ng/ul	92
12) Phenanthrene	17.14	178	4061	0.049	ng/ul#	87
13) Anthracene	17.22	178	3221	0.045	ng/ul#	88
15) Fluoranthene	19.16	202	4614	0.046	ng/ul	84
17) Pyrene	19.52	202	4753	0.049	ng/ul#	84
18) Benzo(a)anthracene	21.24	228	4062	0.046	ng/ul#	83
19) Chrysene	21.30	228	4616	0.050	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	3981	0.042	ng/ul	93
22) Benzo(k)fluoranthene	22.98	252	4839	0.048	ng/ul#	91
23) Benzo(a)pyrene	23.56	252	4146	0.046	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	26.19	276	4595	0.044	ng/ul#	85
25) Dibenzo(a,h)anthracene	26.21	278	3970	0.045	ng/ul#	80
26) Benzo(g,h,i)perylene	26.96	276	4169	0.046	ng/ul	94

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

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