

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081417\
 Data File : BE093793.D
 Acq On : 14 Aug 2017 18:51
 Operator : SJ/JU
 Sample : I4557-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AC4

Quant Time: Aug 14 19:27:36 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 14 18:19:53 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2732	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	14029	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	10842	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	19320	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	19971	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	20582	0.40	ng/ul	0.01

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	5778	0.25	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	29891	0.49	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.75	128	144970	4.248	ng/ul#	93
5) 2-Methylnaphthalene	12.36	142	274063	11.010	ng/ul	99
7) Acenaphthylene	14.24	152	13028	0.287	ng/ul#	47
8) Acenaphthene	14.59	153	27691	0.776	ng/ul	96
9) Fluorene	15.57	166	156861	3.478	ng/ul	95
12) Phenanthrene	17.29	178	182133	3.462	ng/ul#	91
13) Anthracene	17.29	178	178430	3.482	ng/ul	93
15) Fluoranthene	19.30	202	48501	0.737	ng/ul	82
17) Pyrene	19.67	202	74776	1.283	ng/ul#	85
18) Benzo(a)anthracene	21.39	228	19857	0.343	ng/ul#	48
19) Chrysene	21.45	228	17478	0.315	ng/ul#	45
21) Benzo(b)fluoranthene	23.15	252	10691	0.168	ng/ul#	47
22) Benzo(k)fluoranthene	23.15	252	10691	0.170	ng/ul#	44
23) Benzo(a)pyrene	23.80	252	10087	0.167	ng/ul#	75
24) Indeno(1,2,3-cd)pyrene	26.56	276	3681	0.055	ng/ul#	85
26) Benzo(g,h,i)perylene	27.38	276	3277	0.058	ng/ul#	47

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

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