

Data Path : Z:\HPCHEM1\BNA_E\Data\BE061417\
 Data File : BE093197.D
 Acq On : 15 Jun 2017 8:02
 Operator : SJ/MA
 Sample : I3521-18
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5D54

Quant Time: Jun 15 11:14:01 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 15 07:58:47 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	3168	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	17059	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	10896	0.40	ng/ul	0.00
10) Phenanthrene-d10	0.00	188	0	0.00	ng/ul	-17.10
16) Chrysene-d12	21.26	240	24613	0.40	ng/ul	0.00
20) Perylene-d12	23.52	264	1351626	0.40	ng/ul	-0.14
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	7539	0.30	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	20257	0.00	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.57	128	7627	0.180	ng/ul#	93
5) 2-Methylnaphthalene	12.19	142	2259	0.078	ng/ul	98
7) Acenaphthylene	14.09	152	31423	0.672	ng/ul	96
8) Acenaphthene	14.43	153	1393	0.037	ng/ul	95
9) Fluorene	15.42	166	1376	0.031	ng/ul#	82
17) Pyrene	19.52	202	359799	4.957	ng/ul#	87
18) Benzo(a)anthracene	21.24	228	178796	2.723	ng/ul#	85
19) Chrysene	21.29	228	199428	2.928	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	400451	0.098	ng/ul	87
22) Benzo(k)fluoranthene	22.93	252	522022	0.120	ng/ul#	89
23) Benzo(a)pyrene	23.46	252	192496	0.050	ng/ul#	85
26) Benzo(g,h,i)perylene	26.96	276	120881	0.031	ng/ul	97

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

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