

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062017\
 Data File : BE093291.D
 Acq On : 20 Jun 2017 21:00
 Operator : SJ/JU
 Sample : I3627-02
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A08

Quant Time: Jun 21 03:24:01 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 21 03:23:49 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	3	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	127	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	528	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	174	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	759366	0.40	ng/ul	-0.15

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	3797	20.42	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	10668	0.00	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.56	128	4376	13.887	ng/ul#	93
5) 2-Methylnaphthalene	12.18	142	1081	5.017	ng/ul	99
7) Acenaphthylene	14.09	152	13322	5.879	ng/ul	95
8) Acenaphthene	14.42	153	3169	1.734	ng/ul	97
9) Fluorene	15.42	166	5518	2.575	ng/ul	95
17) Pyrene	19.52	202	190478	371.205	ng/ul#	83
18) Benzo(a)anthracene	21.24	228	97928	210.993	ng/ul#	85
19) Chrysene	21.30	228	123514	256.542	ng/ul	98
21) Benzo(b)fluoranthene	22.93	252	180566	0.078	ng/ul	87
22) Benzo(k)fluoranthene	22.93	252	233042	0.095	ng/ul#	89
23) Benzo(a)pyrene	23.45	252	84120	0.039	ng/ul#	85
26) Benzo(g,h,i)perylene	26.96	276	42786	0.020	ng/ul	96

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

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