

Data Path : Z:\HPCHEM1\BNA_E\Data\BE061517\
 Data File : BE093233.D
 Acq On : 16 Jun 2017 8:33
 Operator : SJ/MA
 Sample : I3558-07DL 5X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5C61DL

Quant Time: Jun 16 14:17:49 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 16 05:36:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	2518	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	12994	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	7403	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	18214	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	17216	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	16991	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	592	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	1499	0.03	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.57	128	10199	0.316	ng/ul#	92
5) 2-Methylnaphthalene	12.19	142	1962	0.089	ng/ul	98
7) Acenaphthylene	14.09	152	2143	0.067	ng/ul	92
8) Acenaphthene	14.43	153	2752	0.107	ng/ul	98
9) Fluorene	15.42	166	1800	0.060	ng/ul	92
12) Phenanthrene	17.14	178	14114	0.282	ng/ul#	86
13) Anthracene	17.22	178	5718	0.131	ng/ul#	89
15) Fluoranthene	19.16	202	20722	0.337	ng/ul	84
17) Pyrene	19.52	202	18724	0.369	ng/ul#	85
18) Benzo(a)anthracene	21.24	228	10667	0.232	ng/ul#	86
19) Chrysene	21.29	228	14638	0.307	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	48246	0.936	ng/ul	88
22) Benzo(k)fluoranthene	22.93	252	48246	0.881	ng/ul#	90
23) Benzo(a)pyrene	23.55	252	18137	0.372	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	26.18	276	19549	0.346	ng/ul#	88
25) Dibenzo(a,h)anthracene	26.20	278	4815	0.101	ng/ul#	87
26) Benzo(g,h,i)perylene	26.96	276	17510	0.362	ng/ul	96

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

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