

Data Path : Z:\HPCHEM1\BNA_E\Data\BE101817\
 Data File : BE094438.D
 Acq On : 18 Oct 2017 10:11
 Operator : SJ/JU
 Sample : I5548-09DL 10X
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AC2DL

Quant Time: Oct 18 17:58:51 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 18 17:53:25 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1234	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	6505	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	4024	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	9328	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	9842	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	9630	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.30	152	226	0.02	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	652	0.02	ng/ul	0.00

Target Compounds				Qvalue		
7) Acenaphthylene	14.24	152	764	0.049	ng/ul#	82
9) Fluorene	15.58	166	461	0.030	ng/ul#	89
12) Phenanthrene	17.39	178	2641	0.108	ng/ul	96
13) Anthracene	17.39	178	2985	0.126	ng/ul	98
15) Fluoranthene	19.31	202	31515	0.984	ng/ul	84
17) Pyrene	19.67	202	27715	1.012	ng/ul#	81
18) Benzo(a)anthracene	21.40	228	15560	0.596	ng/ul#	87
19) Chrysene	21.46	228	14211	0.471	ng/ul	97
21) Benzo(b)fluoranthene	23.17	252	24062	0.920	ng/ul	90
22) Benzo(k)fluoranthene	23.17	252	24054	0.795	ng/ul	92
23) Benzo(a)pyrene	23.83	252	11850	0.436	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	26.63	276	7787	0.280	ng/ul#	77
25) Dibenzo(a,h)anthracene	26.62	278	2039	0.087	ng/ul#	67
26) Benzo(g,h,i)perylene	27.47	276	5759	0.240	ng/ul#	89

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

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