

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102717\
 Data File : BE094611.D
 Acq On : 28 Oct 2017 00:23
 Operator : SJ/JU
 Sample : I5842-14DL 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AB2DL

Quant Time: Oct 30 01:12:56 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 30 01:11:00 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1253	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	7371	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	4094	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	8940	0.40	ng/ul	-0.02
16) Chrysene-d12	21.43	240	8886	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	9833	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	833	0.07	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	2043	0.08	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.29	178	702	0.030	ng/ul#	81
15) Fluoranthene	19.31	202	2157	0.079	ng/ul	85
17) Pyrene	19.67	202	2155	0.078	ng/ul	94
18) Benzo(a)anthracene	21.41	228	799	0.031	ng/ul#	28
19) Chrysene	21.47	228	1313	0.050	ng/ul#	51
21) Benzo(b)fluoranthene	23.19	252	2292	0.083	ng/ul#	1
22) Benzo(k)fluoranthene	23.19	252	2303	0.078	ng/ul#	1
23) Benzo(a)pyrene	23.85	252	944	0.034	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	26.68	276	668	0.023	ng/ul#	72
26) Benzo(g,h,i)perylene	27.53	276	602	0.026	ng/ul#	1

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

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