

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102717\
 Data File : BE094634.D
 Acq On : 28 Oct 2017 22:00
 Operator : SJ/JU
 Sample : I5842-09DL 5X
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AA7DL

Quant Time: Oct 30 04:37:58 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 04:29:38 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1949	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	11121	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	6147	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	13263	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	13078	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	13239	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	979	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	2314	0.06	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.29	178	883	0.025	ng/ul#	72
15) Fluoranthene	19.31	202	2051	0.051	ng/ul	83
17) Pyrene	19.67	202	2011	0.050	ng/ul	99
18) Benzo(a)anthracene	21.41	228	1165	0.031	ng/ul#	1
19) Chrysene	21.47	228	1350	0.035	ng/ul#	22
21) Benzo(b)fluoranthene	23.19	252	2406	0.065	ng/ul#	1
22) Benzo(k)fluoranthene	23.19	252	2406	0.060	ng/ul#	1
23) Benzo(a)pyrene	23.85	252	945	0.025	ng/ul#	1
26) Benzo(g,h,i)perylene	27.52	276	679	0.022	ng/ul#	1

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

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