

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102717\
 Data File : BE094636.D
 Acq On : 28 Oct 2017 23:15
 Operator : SJ/JU
 Sample : I5842-10DL 5X
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AA8DL

Quant Time: Oct 30 04:38:18 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	7.91	152	1938	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	10917	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	5874	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	13016	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	12646	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	12607	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.28	152	549	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	1530	0.04	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
12) Phenanthrene	17.29	178	921	0.027	ng/ul#	63
15) Fluoranthene	19.32	202	1846	0.046	ng/ul	83
17) Pyrene	19.68	202	1907	0.049	ng/ul	97
18) Benzo(a)anthracene	21.41	228	897	0.025	ng/ul#	1
19) Chrysene	21.47	228	1321	0.035	ng/ul#	30
21) Benzo(b)fluoranthene	23.19	252	2304	0.065	ng/ul#	1
22) Benzo(k)fluoranthene	23.19	252	2304	0.061	ng/ul#	1
23) Benzo(a)pyrene	23.85	252	923	0.026	ng/ul#	1
26) Benzo(g,h,i)perylene	27.52	276	672	0.023	ng/ul#	1

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

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