

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102717\
 Data File : BE094621.D
 Acq On : 28 Oct 2017 14:03
 Operator : SJ/JU
 Sample : I5842-04DL 5X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AA4DL

Quant Time: Oct 30 02:58:09 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 02:53:50 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1958	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.69	136	11281	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	6228	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	13494	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	13225	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	13926	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	983	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	2418	0.06	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.29	178	3108	0.087	ng/ul#	91
13) Anthracene	17.29	178	3069	0.087	ng/ul	93
15) Fluoranthene	19.31	202	7405	0.180	ng/ul	84
17) Pyrene	19.67	202	6742	0.165	ng/ul#	85
18) Benzo(a)anthracene	21.41	228	2889	0.076	ng/ul#	77
19) Chrysene	21.47	228	3869	0.099	ng/ul#	76
21) Benzo(b)fluoranthene	23.19	252	6799	0.174	ng/ul#	65
22) Benzo(k)fluoranthene	23.19	252	6720	0.160	ng/ul#	62
23) Benzo(a)pyrene	23.85	252	2796	0.071	ng/ul#	55
24) Indeno(1,2,3-cd)pyrene	26.67	276	2099	0.052	ng/ul#	73
26) Benzo(g,h,i)perylene	27.51	276	1666	0.051	ng/ul#	22

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

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