

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102717\
 Data File : BE094627.D
 Acq On : 28 Oct 2017 17:44
 Operator : SJ/JU
 Sample : I5842-19DL 5X
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AB8DL

Quant Time: Oct 30 02:59:10 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	2063	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.69	136	11729	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	6420	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	14174	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	13893	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	14475	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	942	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	2327	0.06	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.29	178	1340	0.036	ng/ul#	92
15) Fluoranthene	19.31	202	3874	0.090	ng/ul	85
17) Pyrene	19.67	202	4842	0.113	ng/ul#	87
18) Benzo(a)anthracene	21.41	228	1883	0.047	ng/ul#	65
19) Chrysene	21.47	228	2888	0.070	ng/ul#	67
21) Benzo(b)fluoranthene	23.19	252	6117	0.150	ng/ul#	72
22) Benzo(k)fluoranthene	23.19	252	6085	0.139	ng/ul#	69
23) Benzo(a)pyrene	23.85	252	2265	0.056	ng/ul#	52
24) Indeno(1,2,3-cd)pyrene	26.67	276	1608	0.038	ng/ul#	76
26) Benzo(g,h,i)perylene	27.51	276	1199	0.035	ng/ul#	27

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

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