

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102717\
 Data File : BE094625.D
 Acq On : 28 Oct 2017 16:29
 Operator : SJ/JU
 Sample : I5842-02DL 5X
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AA2DL

Quant Time: Oct 30 02:58:52 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	2008	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.69	136	11524	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	6398	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	14200	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	14076	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	14786	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	977	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	2401	0.06	ng/ul	0.00
Target Compounds				Ovalue		
11) Pentachlorophenol	16.98	266	15	0.021	ng/ul#	20
12) Phenanthrene	17.29	178	2167	0.058	ng/ul#	93
15) Fluoranthene	19.31	202	7195	0.166	ng/ul	83
17) Pyrene	19.67	202	6288	0.145	ng/ul#	86
18) Benzo(a)anthracene	21.41	228	2704	0.067	ng/ul#	59
19) Chrysene	21.46	228	4267	0.103	ng/ul#	64
21) Benzo(b)fluoranthene	23.18	252	7870	0.189	ng/ul#	59
22) Benzo(k)fluoranthene	23.18	252	7818	0.175	ng/ul#	56
23) Benzo(a)pyrene	23.85	252	3014	0.072	ng/ul#	41
24) Indeno(1,2,3-cd)pyrene	26.66	276	2101	0.049	ng/ul#	76
26) Benzo(g,h,i)perylene	27.51	276	1887	0.055	ng/ul#	26

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

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