

Data Path : Z:\HPCHEM1\BNA E\DATA\BE103117\
 Data File : BE094657.D
 Acq On : 31 Oct 2017 14:22
 Operator : SJ/JU
 Sample : I5842-03DL 10X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AA3DL

Manual Integrations
 APPROVED

Sohil
 11/1/2017 4:29:40 PM

Quant Time: Nov 01 02:11:59 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 01 01:43:26 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1543	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	9412	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	5344	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	11755	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	12030	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	11819	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	327	0.02	ng/ul	0.01
14) Fluoranthene-d10	19.28	212	744	0.02	ng/ul	0.00

Target Compounds						Ovalue
8) Acenaphthene	14.58	153	564	0.032	ng/ul#	93
9) Fluorene	15.57	166	485	0.025	ng/ul#	86
12) Phenanthrene	17.29	178	9976	0.321	ng/ul#	88
13) Anthracene	17.40	178	1919	0.063	ng/ul	94
15) Fluoranthene	19.31	202	28246	0.787	ng/ul	84
17) Pyrene	19.67	202	22284	0.599	ng/ul#	81
18) Benzo(a)anthracene	21.41	228	6822	0.198	ng/ul#	88
19) Chrysene	21.46	228	12056	0.339	ng/ul	98
21) Benzo(b)fluoranthene	23.19	252	9532m	0.287	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	3315m	0.093	ng/ul	
23) Benzo(a)pyrene	23.85	252	4135	0.124	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	26.67	276	2309	0.067	ng/ul#	74
26) Benzo(g,h,i)perylene	27.51	276	1732	0.063	ng/ul#	64

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

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