

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110317\
 Data File : BE094725.D
 Acq On : 4 Nov 2017 1:38
 Operator : SJ/JU
 Sample : I5955-13 5X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-6(3.5-4.5)-101717

Manual Integrations
 APPROVED

Sohil
 11/4/2017 2:16:54 PM

Quant Time: Nov 04 01:39:41 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 04 02:38:13 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1503	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.69	136	8975	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	4965	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	10773	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	10704	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	11635	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	618	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	1515	0.05	ng/ul	0.00
Target Compounds				Qvalue		
12) Phenanthrene	17.29	178	2782	0.098	ng/ul#	91
15) Fluoranthene	19.31	202	4656	0.142	ng/ul	84
17) Pyrene	19.67	202	4188	0.127	ng/ul#	84
18) Benzo(a)anthracene	21.41	228	2124	0.069	ng/ul#	73
19) Chrysene	21.47	228	2423	0.077	ng/ul#	68
21) Benzo(b)fluoranthene	23.19	252	2554m	0.078	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	886m	0.025	ng/ul	
23) Benzo(a)pyrene	23.85	252	1601	0.049	ng/ul#	23
24) Indeno(1,2,3-cd)pyrene	26.68	276	819	0.024	ng/ul#	73
26) Benzo(g,h,i)perylene	27.53	276	805	0.030	ng/ul#	8

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

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