

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110317\
 Data File : BE094723.D
 Acq On : 4 Nov 2017 00:29
 Operator : SJ/JU
 Sample : I5955-10 5X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 DUP-6

Manual Integrations
 APPROVED

Sohil
 11/4/2017 2:16:52 PM

Quant Time: Nov 04 00:43:47 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 01:15:30 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1610	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.69	136	9768	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.51	164	5516	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	11617	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	11644	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	13394	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	1114	0.07	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	3358	0.10	ng/ul	0.00
Target Compounds				Qvalue		
12) Phenanthrene	17.29	178	3717	0.121	ng/ul	94
13) Anthracene	17.38	178	1136m	0.038	ng/ul	
15) Fluoranthene	19.31	202	17017	0.480	ng/ul	84
17) Pyrene	19.67	202	19471	0.541	ng/ul#	80
18) Benzo(a)anthracene	21.40	228	13748	0.413	ng/ul#	88
19) Chrysene	21.46	228	12531	0.364	ng/ul#	87
21) Benzo(b)fluoranthene	23.18	252	16348m	0.434	ng/ul	
22) Benzo(k)fluoranthene	23.22	252	6500m	0.161	ng/ul	
23) Benzo(a)pyrene	23.84	252	12265	0.326	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	26.66	276	6814	0.175	ng/ul#	71
25) Dibenzo(a,h)anthracene	26.65	278	1605	0.049	ng/ul#	1
26) Benzo(g,h,i)perylene	27.50	276	6283	0.201	ng/ul#	70

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

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