

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110317\
 Data File : BE094717.D
 Acq On : 3 Nov 2017 20:54
 Operator : SJ/JU
 Sample : I5902-02 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-5(5-6.5)-101317

Manual Integrations
 APPROVED

Sohil
 11/4/2017 2:15:39 PM

Quant Time: Nov 04 00:27:14 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	1716	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.69	136	10343	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	5686	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	12045	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	12195	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	13362	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	679	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	1626	0.05	ng/ul	0.00
Target Compounds				Qvalue		
12) Phenanthrene	17.29	178	1893	0.060	ng/ul#	90
15) Fluoranthene	19.31	202	3895	0.106	ng/ul	84
17) Pyrene	19.67	202	4012	0.106	ng/ul#	87
18) Benzo(a)anthracene	21.41	228	2284	0.066	ng/ul#	75
19) Chrysene	21.47	228	2176	0.060	ng/ul#	74
21) Benzo(b)fluoranthene	23.18	252	2574m	0.068	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	1191m	0.030	ng/ul	
23) Benzo(a)pyrene	23.84	252	1878	0.050	ng/ul#	48
24) Indeno(1,2,3-cd)pyrene	26.67	276	1052	0.027	ng/ul#	76
26) Benzo(g,h,i)perylene	27.52	276	916	0.029	ng/ul#	23

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

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