

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101417\
 Data File : BE094379.D
 Acq On : 15 Oct 2017 5:49
 Operator : SJ/JU
 Sample : I5548-15
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AF5

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:25:45 PM

Quant Time: Oct 16 06:21:03 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 05:09:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1570	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	8444	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	5218	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	11468	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	10905	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	10887m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	2989	0.23	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	8040	0.21	ng/ul	0.00
Target Compounds					Ovalue	
12) Phenanthrene	17.29	178	7128	0.238	ng/ul#	88
13) Anthracene	17.38	178	1795	0.062	ng/ul#	80
15) Fluoranthene	19.31	202	13277	0.337	ng/ul	84
17) Pyrene	19.67	202	12579	0.415	ng/ul#	84
18) Benzo(a)anthracene	21.40	228	6001	0.207	ng/ul#	81
19) Chrysene	21.46	228	5276	0.158	ng/ul#	73
21) Benzo(b)fluoranthene	23.17	252	5793	0.196	ng/ul#	71
22) Benzo(k)fluoranthene	23.21	252	2732m	0.080	ng/ul	
23) Benzo(a)pyrene	23.83	252	4936m	0.161	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.62	276	2974m	0.095	ng/ul	
26) Benzo(g,h,i)perylene	27.46	276	2999m	0.111	ng/ul	

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

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