

Data Path : Z:\HPCHEM1\BNA E\DATA\BE100917\
 Data File : BE094223.D
 Acq On : 9 Oct 2017 16:28
 Operator : SJ/JU
 Sample : I5522-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AD2

Manual Integrations
 APPROVED

Sohil
 10/11/2017 9:29:45 AM

Quant Time: Oct 10 16:52:34 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 10 09:18:40 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1269	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	7042	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	4259	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	9578m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	10058	0.40	ng/ul	0.00
20) Perylene-d12	23.93	264	8452m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	3110	0.28	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	8673	0.27	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.29	178	4220	0.169	ng/ul	92
15) Fluoranthene	19.30	202	12806	0.389	ng/ul	84
17) Pyrene	19.66	202	12087	0.432	ng/ul#	85
18) Benzo(a)anthracene	21.39	228	7022	0.263	ng/ul#	89
19) Chrysene	21.45	228	8912	0.289	ng/ul#	88
21) Benzo(b)fluoranthene	23.15	252	9308m	0.405	ng/ul	
22) Benzo(k)fluoranthene	23.20	252	3450m	0.130	ng/ul	
23) Benzo(a)pyrene	23.81	252	6585m	0.276	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.57	276	5884m	0.241	ng/ul	
26) Benzo(g,h,i)perylene	27.41	276	5622m	0.267	ng/ul	

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

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