

Data Path : Z:\HPCHEM1\BNA E\DATA\BE100917\
 Data File : BE094229.D
 Acq On : 9 Oct 2017 20:12
 Operator : SJ/JU
 Sample : I5522-10
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AD9

Manual Integrations
 APPROVED

Sohil
 10/11/2017 9:24:18 AM

Quant Time: Oct 10 17:20:35 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	2446	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	12938	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	7498	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.24	188	16708	0.40	ng/ul	-0.02
16) Chrysene-d12	21.41	240	16717	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	15401m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	1184	0.06	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	3385	0.06	ng/ul	0.00
Target Compounds						
12) Phenanthrene	17.29	178	1950	0.045	ng/ul#	96
15) Fluoranthene	19.30	202	5370	0.094	ng/ul	85
17) Pyrene	19.66	202	5549	0.119	ng/ul#	88
18) Benzo(a)anthracene	21.39	228	4799	0.108	ng/ul#	88
19) Chrysene	21.44	228	6005	0.117	ng/ul#	87
21) Benzo(b)fluoranthene	23.15	252	11327m	0.271	ng/ul	
22) Benzo(k)fluoranthene	23.19	252	4148m	0.086	ng/ul	
23) Benzo(a)pyrene	23.80	252	7912m	0.182	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.57	276	5942m	0.134	ng/ul	
26) Benzo(g,h,i)perylene	27.40	276	5087m	0.133	ng/ul	

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

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