

Data Path : Z:\HPCHEM1\BNA E\DATA\BE100917\
 Data File : BE094228.D
 Acq On : 9 Oct 2017 19:36
 Operator : SJ/JU
 Sample : I5522-09
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AD8

Manual Integrations
 APPROVED

Sohil
 10/11/2017 9:24:16 AM

Quant Time: Oct 10 17:15:05 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	2437	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	12863	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	7487	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.24	188	16816	0.40	ng/ul	-0.02
16) Chrysene-d12	21.41	240	16464	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	15260m	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	1636	0.08	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	8837	0.16	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.29	178	3260	0.074	ng/ul	96
15) Fluoranthene	19.30	202	12293	0.213	ng/ul	84
17) Pyrene	19.66	202	12327	0.269	ng/ul#	86
18) Benzo(a)anthracene	21.39	228	6108	0.140	ng/ul#	90
19) Chrysene	21.45	228	8212	0.163	ng/ul#	90
21) Benzo(b)fluoranthene	23.15	252	11752m	0.283	ng/ul	
22) Benzo(k)fluoranthene	23.19	252	3671m	0.077	ng/ul	
23) Benzo(a)pyrene	23.81	252	6918m	0.161	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.57	276	5129m	0.116	ng/ul	
26) Benzo(g,h,i)perylene	27.40	276	5102m	0.134	ng/ul	

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

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