

Data Path : Z:\HPCHEM1\BNA E\DATA\BE101017\
 Data File : BE094251.D
 Acq On : 10 Oct 2017 13:03
 Operator : SJ/JU
 Sample : I5522-04
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AD5

Manual Integrations
 APPROVED

Sohil
 10/11/2017 7:19:39 PM

Quant Time: Oct 11 11:23:04 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 11 08:24:24 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1723	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	9092	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	5441	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.24	188	11635m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.41	240	11470	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	10216m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	3734	0.26	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	9208	0.24	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	2361	0.106	ng/ul#	94
5) 2-Methylnaphthalene	12.35	142	491	0.034	ng/ul	99
12) Phenanthrene	17.29	178	6532	0.215	ng/ul#	93
15) Fluoranthene	19.30	202	18891	0.473	ng/ul	83
17) Pyrene	19.66	202	17936	0.562	ng/ul#	83
18) Benzo(a)anthracene	21.39	228	9028	0.297	ng/ul#	90
19) Chrysene	21.45	228	10309	0.293	ng/ul	93
21) Benzo(b)fluoranthene	23.15	252	13658m	0.492	ng/ul	
22) Benzo(k)fluoranthene	23.19	252	4460m	0.139	ng/ul	
23) Benzo(a)pyrene	23.81	252	9579m	0.332	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.58	276	6030m	0.204	ng/ul	
26) Benzo(g,h,i)perylene	27.40	276	5102m	0.201	ng/ul	

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

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