

Data Path : Z:\HPCHEM1\BNA E\DATA\BE101017\
 Data File : BE094257.D
 Acq On : 10 Oct 2017 18:48
 Operator : SJ/JU
 Sample : I5522-11
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AE0

Manual Integrations
 APPROVED

Sohil
 10/11/2017 7:19:51 PM

Quant Time: Oct 11 07:38:28 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 07:23:41 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1964	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	10215	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	5680	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.24	188	11807	0.40	ng/ul	-0.02
16) Chrysene-d12	21.41	240	12158	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	13438	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	5289	0.33	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	12076	0.31	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.29	178	2185	0.071	ng/ul	96
15) Fluoranthene	19.30	202	5900	0.146	ng/ul	85
17) Pyrene	19.66	202	5944	0.176	ng/ul#	87
18) Benzo(a)anthracene	21.40	228	2511	0.078	ng/ul#	87
19) Chrysene	21.45	228	3494	0.094	ng/ul#	87
21) Benzo(b)fluoranthene	23.15	252	4693m	0.129	ng/ul	
22) Benzo(k)fluoranthene	23.19	252	1671m	0.040	ng/ul	
23) Benzo(a)pyrene	23.81	252	2651	0.070	ng/ul#	77
24) Indeno(1,2,3-cd)pyrene	26.58	276	2206	0.057	ng/ul#	81
26) Benzo(g,h,i)perylene	27.40	276	1877	0.056	ng/ul#	62

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

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