

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101117\
 Data File : BE094278.D
 Acq On : 11 Oct 2017 13:08
 Operator : SJ/JU
 Sample : I5522-10
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AD9

Manual Integrations
 APPROVED

Sohil
 10/13/2017 12:05:27 PM

Quant Time: Oct 12 07:29:18 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 12 07:23:36 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	2067	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	10848	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	6324	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.24	188	13455	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	13634	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	13704m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	4202	0.25	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	10144	0.23	ng/ul	0.00
Target Compounds					Ovalue	
12) Phenanthrene	17.29	178	2075	0.059	ng/ul	97
15) Fluoranthene	19.30	202	5916	0.128	ng/ul	85
17) Pyrene	19.66	202	6112	0.161	ng/ul#	87
18) Benzo(a)anthracene	21.39	228	5253	0.145	ng/ul#	88
19) Chrysene	21.44	228	6502	0.156	ng/ul	93
21) Benzo(b)fluoranthene	23.15	252	11503	0.309	ng/ul	93
22) Benzo(k)fluoranthene	23.19	252	4382m	0.102	ng/ul	
23) Benzo(a)pyrene	23.80	252	8525m	0.220	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.57	276	6131m	0.155	ng/ul	
25) Dibenzo(a,h)anthracene	26.58	278	1894m	0.057	ng/ul	
26) Benzo(g,h,i)perylene	27.40	276	5125m	0.150	ng/ul	

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

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