

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101417\
 Data File : BE094358.D
 Acq On : 14 Oct 2017 16:56
 Operator : SJ/JU
 Sample : I5522-13
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AE2

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:25:04 PM

Quant Time: Oct 15 01:41:58 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 14 05:22:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1175	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.70	136	5884	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	3654	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	9514	0.40	ng/ul	-0.02
16) Chrysene-d12	21.42	240	10249	0.40	ng/ul	-0.01
20) Perylene-d12	23.94	264	9370m	0.40	ng/ul	-0.03
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	317	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	2344	0.07	ng/ul	-0.01
Target Compounds				Ovalue		
12) Phenanthrene	17.29	178	885	0.036	ng/ul#	85
15) Fluoranthene	19.31	202	3014	0.092	ng/ul	85
17) Pyrene	19.67	202	3196	0.112	ng/ul#	86
18) Benzo(a)anthracene	21.41	228	2173	0.080	ng/ul#	77
19) Chrysene	21.46	228	3089	0.098	ng/ul#	78
21) Benzo(b)fluoranthene	23.17	252	4461m	0.175	ng/ul	
22) Benzo(k)fluoranthene	23.21	252	1341m	0.046	ng/ul	
23) Benzo(a)pyrene	23.83	252	2660m	0.101	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.63	276	2139m	0.079	ng/ul	
26) Benzo(g,h,i)perylene	27.47	276	2084m	0.089	ng/ul	

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

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