

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102717\
 Data File : BE094637.D
 Acq On : 28 Oct 2017 23:51
 Operator : SJ/JU
 Sample : I5842-11
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AA9

Manual Integrations
 APPROVED

Sohil
 10/30/2017 6:40:25 PM

Quant Time: Oct 30 05:00:08 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 30 04:29:38 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1924	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	11076	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	6278	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	13304m	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	12604	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	12428m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	5348	0.30	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	12873	0.33	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.29	178	2444m	0.070	ng/ul	
15) Fluoranthene	19.32	202	8713	0.214	ng/ul	83
17) Pyrene	19.68	202	8572	0.220	ng/ul#	85
18) Benzo(a)anthracene	21.41	228	3167	0.088	ng/ul#	83
19) Chrysene	21.47	228	4051	0.109	ng/ul#	83
21) Benzo(b)fluoranthene	23.19	252	4861m	0.139	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	1775m	0.047	ng/ul	
23) Benzo(a)pyrene	23.85	252	2517m	0.072	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.68	276	1654m	0.046	ng/ul	
26) Benzo(g,h,i)perylene	27.52	276	1243m	0.043	ng/ul	

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

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