

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102717\
 Data File : BE094626.D
 Acq On : 28 Oct 2017 17:08
 Operator : SJ/JU
 Sample : I5842-19
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AB8

Manual Integrations
 APPROVED

Sohil
 10/30/2017 6:40:08 PM

Quant Time: Oct 30 04:10:17 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 02:53:50 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1552	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.69	136	8823	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	5029	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	10838	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	10300	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	10793m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	3452	0.24	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	8688	0.27	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.29	178	4836	0.169	ng/ul#	91
13) Anthracene	17.38	178	1100	0.039	ng/ul#	70
15) Fluoranthene	19.31	202	14329	0.433	ng/ul	84
17) Pyrene	19.67	202	17445	0.548	ng/ul#	82
18) Benzo(a)anthracene	21.41	228	6902	0.234	ng/ul#	89
19) Chrysene	21.46	228	10229	0.336	ng/ul#	87
21) Benzo(b)fluoranthene	23.19	252	15960m	0.526	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	4991m	0.153	ng/ul	
23) Benzo(a)pyrene	23.85	252	7909m	0.261	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.67	276	5755m	0.183	ng/ul	
25) Dibenzo(a,h)anthracene	26.66	278	1318m	0.050	ng/ul	
26) Benzo(g,h,i)perylene	27.51	276	4656m	0.185	ng/ul	

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

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