

Data Path : Z:\HPCHEM1\BNA E\DATA\BE101017\
 Data File : BE094258.D
 Acq On : 10 Oct 2017 19:27
 Operator : SJ/JU
 Sample : I5522-18
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AC4

Manual Integrations
 APPROVED

Sohil
 10/11/2017 7:19:53 PM

Quant Time: Oct 11 07:40:31 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	1778	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	9154	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	5215	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.24	188	11221	0.40	ng/ul	-0.02
16) Chrysene-d12	21.41	240	11552	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	12688	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	5034	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	12704	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	2408	0.107	ng/ul#	94
5) 2-Methylnaphthalene	12.35	142	778	0.053	ng/ul	98
7) Acenaphthylene	14.24	152	1173	0.058	ng/ul#	88
8) Acenaphthene	14.57	153	2621	0.157	ng/ul	99
9) Fluorene	15.56	166	2746	0.140	ng/ul	94
12) Phenanthrene	17.27	178	59559	2.031	ng/ul#	83
13) Anthracene	17.38	178	13338	0.468	ng/ul#	92
15) Fluoranthene	19.30	202	108332	2.812	ng/ul	84
17) Pyrene	19.66	202	94213	2.931	ng/ul#	84
18) Benzo(a)anthracene	21.39	228	51716	1.688	ng/ul#	87
19) Chrysene	21.44	228	48878	1.381	ng/ul	99
21) Benzo(b)fluoranthene	23.15	252	55803m	1.619	ng/ul	
22) Benzo(k)fluoranthene	23.19	252	21403m	0.537	ng/ul	
23) Benzo(a)pyrene	23.80	252	42962	1.200	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	26.57	276	25570	0.698	ng/ul#	79
25) Dibenzo(a,h)anthracene	26.57	278	6770	0.220	ng/ul#	94
26) Benzo(a,h,i)perylene	27.39	276	18009	0.570	ng/ul	98

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

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