

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101417\
 Data File : BE094371.D
 Acq On : 15 Oct 2017 00:54
 Operator : SJ/JU
 Sample : I5548-07
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AB8

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:25:30 PM

Quant Time: Oct 16 05:18:08 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 05:09:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1476	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	8142	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	5121	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	11454m	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	9906	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	9353m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	3593	0.28	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	8655	0.23	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.75	128	1654	0.083	ng/ul#	72
5) 2-Methylnaphthalene	12.36	142	1511	0.117	ng/ul	100
7) Acenaphthylene	14.25	152	835	0.042	ng/ul#	67
8) Acenaphthene	14.59	153	1452	0.088	ng/ul	97
9) Fluorene	15.57	166	1333	0.069	ng/ul#	70
12) Phenanthrene	17.29	178	23299	0.778	ng/ul#	89
13) Anthracene	17.40	178	4837	0.166	ng/ul#	78
15) Fluoranthene	19.32	202	51783	1.317	ng/ul	83
17) Pyrene	19.68	202	46878	1.701	ng/ul#	82
18) Benzo(a)anthracene	21.41	228	24360	0.927	ng/ul#	74
19) Chrysene	21.46	228	27015	0.890	ng/ul#	70
21) Benzo(b)fluoranthene	23.18	252	32121	1.264	ng/ul#	67
22) Benzo(k)fluoranthene	23.22	252	12588m	0.429	ng/ul	
23) Benzo(a)pyrene	23.84	252	22045m	0.835	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.64	276	16385m	0.607	ng/ul	
25) Dibenzo(a,h)anthracene	26.64	278	3556m	0.157	ng/ul	
26) Benzo(a,h,i)perylene	27.48	276	12765m	0.548	ng/ul	

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

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