

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101417\
 Data File : BE094402.D
 Acq On : 15 Oct 2017 20:35
 Operator : SJ/JU
 Sample : I5548-13
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AC9

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:26:37 PM

Quant Time: Oct 16 08:11:29 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 07:45:00 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1837	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	9987	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	5781	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	12100	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	11538	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	11735m	0.40	ng/ul	0.03
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	2832	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	6332	0.16	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.75	128	3565	0.146	ng/ul#	95
5) 2-Methylnaphthalene	12.36	142	2407	0.152	ng/ul	98
7) Acenaphthylene	14.25	152	3593	0.159	ng/ul	95
8) Acenaphthene	14.59	153	2354	0.127	ng/ul	99
9) Fluorene	15.57	166	2167	0.100	ng/ul#	93
12) Phenanthrene	17.29	178	107799m	3.409	ng/ul	
13) Anthracene	17.39	178	15210	0.495	ng/ul	95
15) Fluoranthene	19.32	202	551410	13.272	ng/ul	84
17) Pyrene	19.68	202	495675	15.440	ng/ul#	80
18) Benzo(a)anthracene	21.41	228	249449	8.150	ng/ul#	88
19) Chrysene	21.47	228	459822	13.008	ng/ul	98
21) Benzo(b)fluoranthene	23.20	252	653075	20.487	ng/ul	86
22) Benzo(k)fluoranthene	23.24	252	212936m	5.779	ng/ul	
23) Benzo(a)pyrene	23.86	252	381719m	11.523	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.68	276	330926m	9.766	ng/ul	
25) Dibenzo(a,h)anthracene	26.68	278	80753m	2.841	ng/ul	
26) Benzo(a,h,i)perylene	27.53	276	288965m	9.886	ng/ul	

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

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