

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101417\
 Data File : BE094403.D
 Acq On : 15 Oct 2017 21:14
 Operator : SJ/JU
 Sample : I5548-13DL 5X
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AC9DL

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:26:39 PM

Quant Time: Oct 16 08:15:13 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	1945	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	10381	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	5871	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	12290	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	12587	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	11665	0.40	ng/ul	0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	645	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	1432	0.03	ng/ul	0.00
Target Compounds						
					Ovalue	
3) Naphthalene	10.76	128	870	0.034	ng/ul#	80
5) 2-Methylnaphthalene	12.36	142	535	0.032	ng/ul	100
7) Acenaphthylene	14.25	152	840	0.037	ng/ul#	75
8) Acenaphthene	14.59	153	547	0.029	ng/ul	95
9) Fluorene	15.57	166	543	0.025	ng/ul#	78
12) Phenanthrene	17.31	178	24703	0.769	ng/ul	95
13) Anthracene	17.40	178	3506m	0.112	ng/ul	
15) Fluoranthene	19.32	202	128687	3.050	ng/ul	84
17) Pyrene	19.68	202	113821	3.250	ng/ul#	79
18) Benzo(a)anthracene	21.41	228	60658	1.817	ng/ul#	88
19) Chrysene	21.47	228	105696	2.741	ng/ul	95
21) Benzo(b)fluoranthene	23.19	252	165849m	5.234	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	57916m	1.581	ng/ul	
23) Benzo(a)pyrene	23.86	252	90674	2.754	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	26.66	276	81388	2.416	ng/ul#	76
25) Dibenzo(a,h)anthracene	26.66	278	19215	0.680	ng/ul#	71
26) Benzo(a,h,i)perylene	27.50	276	72607	2.499	ng/ul	94

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

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