

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101417\
 Data File : BE094406.D
 Acq On : 15 Oct 2017 23:04
 Operator : SJ/JU
 Sample : I5548-20
 Misc :
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AD1

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:26:45 PM

Quant Time: Oct 16 08:23:44 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	1968	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	10496	0.40	ng/ul	0.01
6) Acenaphthene-d10	14.53	164	6111	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	12235	0.40	ng/ul	0.02
16) Chrysene-d12	21.44	240	11426	0.40	ng/ul	0.02
20) Perylene-d12	23.99	264	12945m	0.40	ng/ul	0.04
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	3233	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.30	212	7040m	0.17	ng/ul	0.01
Target Compounds				Ovalue		
3) Naphthalene	10.76	128	11474	0.446	ng/ul#	92
5) 2-Methylnaphthalene	12.36	142	6889	0.413	ng/ul	99
7) Acenaphthylene	14.26	152	15256	0.640	ng/ul	98
8) Acenaphthene	14.59	153	16437	0.838	ng/ul	97
9) Fluorene	15.58	166	18630	0.811	ng/ul	95
12) Phenanthrene	17.31	178	491575	15.375	ng/ul#	89
13) Anthracene	17.40	178	151869	4.884	ng/ul#	89
15) Fluoranthene	19.33	202	1636301	38.950	ng/ul	84
17) Pyrene	19.69	202	1635416	51.443	ng/ul#	84
18) Benzo(a)anthracene	21.43	228	929322	30.659	ng/ul#	86
19) Chrysene	21.48	228	820527	23.440	ng/ul	98
21) Benzo(b)fluoranthene	23.21	252	1000330	28.447	ng/ul	86
22) Benzo(k)fluoranthene	23.25	252	361210m	8.887	ng/ul	
23) Benzo(a)pyrene	23.88	252	771259m	21.107	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.71	276	464474m	12.426	ng/ul	
25) Dibenzo(a,h)anthracene	26.71	278	123839m	3.950	ng/ul	
26) Benzo(a,h,i)perylene	27.55	276	353647m	10.968	ng/ul	

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

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