

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101417\
 Data File : BE094383.D
 Acq On : 15 Oct 2017 8:15
 Operator : SJ/JU
 Sample : I5548-18DL 5X
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AF8DL

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:25:53 PM

Quant Time: Oct 16 06:39:08 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	2160	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	11268	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	7022	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	13496	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	14004	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	11939	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	737	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	1750	0.04	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	97332	3.527	ng/ul#	92
5) 2-Methylnaphthalene	12.35	142	125891	7.031	ng/ul	98
8) Acenaphthene	14.59	153	4852	0.215	ng/ul#	64
9) Fluorene	15.57	166	1545	0.059	ng/ul#	63
12) Phenanthrene	17.29	178	24718	0.701	ng/ul#	91
13) Anthracene	17.38	178	3138	0.091	ng/ul#	64
15) Fluoranthene	19.31	202	20509	0.443	ng/ul	82
17) Pyrene	19.67	202	20126	0.517	ng/ul#	87
18) Benzo(a)anthracene	21.40	228	10509	0.283	ng/ul#	58
19) Chrysene	21.46	228	11964	0.279	ng/ul#	54
21) Benzo(b)fluoranthene	23.18	252	13546m	0.418	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	4455m	0.119	ng/ul	
23) Benzo(a)pyrene	23.85	252	8595	0.255	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	26.65	276	5545	0.161	ng/ul#	80
26) Benzo(g,h,i)perylene	27.49	276	6189	0.208	ng/ul#	28

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

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