

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101117\
 Data File : BE094286.D
 Acq On : 11 Oct 2017 19:12
 Operator : SJ/JU
 Sample : I5490-01DL 5X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-43(0-1)-092617DL

Manual Integrations
 APPROVED

Sohil
 10/13/2017 12:05:39 PM

Quant Time: Oct 12 08:30:56 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 12 07:31:23 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1870	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	9446	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	5239	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	10975	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	10866	0.40	ng/ul	0.00
20) Perylene-d12	23.93	264	11416	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	659	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	1529	0.04	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	7380	0.319	ng/ul#	94
5) 2-Methylnaphthalene	12.35	142	3660	0.244	ng/ul	97
7) Acenaphthylene	14.24	152	977	0.048	ng/ul#	64
8) Acenaphthene	14.58	153	1586	0.094	ng/ul	99
9) Fluorene	15.57	166	1773	0.090	ng/ul#	89
12) Phenanthrene	17.29	178	23309	0.813	ng/ul#	92
13) Anthracene	17.38	178	5898	0.211	ng/ul#	95
15) Fluoranthene	19.30	202	31596	0.838	ng/ul	84
17) Pyrene	19.67	202	31906	1.055	ng/ul#	82
18) Benzo(a)anthracene	21.40	228	15521	0.538	ng/ul#	79
19) Chrysene	21.45	228	17960	0.540	ng/ul#	74
21) Benzo(b)fluoranthene	23.16	252	18685m	0.603	ng/ul	
22) Benzo(k)fluoranthene	23.20	252	6452m	0.180	ng/ul	
23) Benzo(a)pyrene	23.81	252	13426	0.417	ng/ul#	60
24) Indeno(1,2,3-cd)pyrene	26.59	276	9250	0.281	ng/ul#	76
26) Benzo(a,h,i)perylene	27.42	276	13340	0.469	ng/ul#	65

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

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