

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101117\
 Data File : BE094296.D
 Acq On : 12 Oct 2017 1:22
 Operator : SJ/JU
 Sample : I5490-07
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-41(5-7)-092717

Manual Integrations
 APPROVED

Sohil
 10/13/2017 12:05:56 PM

Quant Time: Oct 12 09:14:48 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 08:55:52 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	1740	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	8619	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	5092	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	10774m	0.40	ng/ul	0.02
16) Chrysene-d12	21.44	240	9796	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	9189m	0.40	ng/ul	0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.30	152	4276	0.32	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	9867	0.27	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.76	128	1473	0.070	ng/ul#	88
5) 2-Methylnaphthalene	12.37	142	506	0.037	ng/ul	97
7) Acenaphthylene	14.26	152	1264	0.064	ng/ul#	78
8) Acenaphthene	14.60	153	559	0.034	ng/ul	93
9) Fluorene	15.58	166	589	0.031	ng/ul#	70
12) Phenanthrene	17.31	178	19556	0.695	ng/ul#	91
13) Anthracene	17.39	178	5611	0.205	ng/ul#	82
15) Fluoranthene	19.33	202	64343	1.739	ng/ul	84
17) Pyrene	19.69	202	70319	2.580	ng/ul#	80
18) Benzo(a)anthracene	21.42	228	48910	1.882	ng/ul#	89
19) Chrysene	21.47	228	47272	1.575	ng/ul#	84
21) Benzo(b)fluoranthene	23.20	252	50623	2.028	ng/ul#	69
22) Benzo(k)fluoranthene	23.24	252	21603m	0.749	ng/ul	
23) Benzo(a)pyrene	23.86	252	41488m	1.599	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.66	276	21566m	0.813	ng/ul	
25) Dibenzo(a,h)anthracene	26.66	278	5926m	0.266	ng/ul	
26) Benzo(a,h,i)perylene	27.51	276	16316m	0.713	ng/ul	

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

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