

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101117\
 Data File : BE094285.D
 Acq On : 11 Oct 2017 18:36
 Operator : SJ/JU
 Sample : I5490-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-70(0-1)-092717

Manual Integrations
 APPROVED

Sohil
 10/13/2017 12:05:37 PM

Quant Time: Oct 12 08:27:44 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : 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	1825	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	9299	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	5430	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	11070	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	11536	0.40	ng/ul	0.00
20) Perylene-d12	23.94	264	10648m	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	2669	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	6263m	0.17	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.74	128	6116	0.269	ng/ul#	96
5) 2-Methylnaphthalene	12.35	142	2870	0.194	ng/ul	99
7) Acenaphthylene	14.24	152	6647	0.314	ng/ul	98
8) Acenaphthene	14.57	153	3618	0.208	ng/ul	97
9) Fluorene	15.57	166	4967	0.243	ng/ul	93
12) Phenanthrene	17.29	178	90931	3.143	ng/ul#	92
13) Anthracene	17.38	178	23764	0.845	ng/ul#	93
15) Fluoranthene	19.30	202	179524	4.723	ng/ul	84
17) Pyrene	19.66	202	165567	5.158	ng/ul#	84
18) Benzo(a)anthracene	21.40	228	89910	2.938	ng/ul#	88
19) Chrysene	21.45	228	88575	2.506	ng/ul	93
21) Benzo(b)fluoranthene	23.16	252	112426	3.887	ng/ul	86
22) Benzo(k)fluoranthene	23.20	252	43113m	1.289	ng/ul	
23) Benzo(a)pyrene	23.82	252	82858m	2.757	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.60	276	61412m	1.997	ng/ul	
25) Dibenzo(a,h)anthracene	26.60	278	18522m	0.718	ng/ul	
26) Benzo(a,h,i)perylene	27.43	276	61206m	2.308	ng/ul	

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

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