

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110617\
 Data File : BE094759.D
 Acq On : 6 Nov 2017 18:39
 Operator : SJ/JU
 Sample : I6096-01 5X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-76(0-1)-102317

Manual Integrations
 APPROVED

Sohil
 11/7/2017 5:41:49 PM

Quant Time: Nov 07 01:17:33 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 07 00:12:32 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1276	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	7512	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	4289	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	9190	0.40	ng/ul	-0.02
16) Chrysene-d12	21.43	240	9145	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	9476	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	375	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	1033	0.04	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.74	128	5826	0.317	ng/ul#	94
5) 2-Methylnaphthalene	12.35	142	2584	0.201	ng/ul	97
7) Acenaphthylene	14.24	152	3880	0.200	ng/ul	97
8) Acenaphthene	14.57	153	8418	0.604	ng/ul	96
9) Fluorene	15.57	166	8407	0.537	ng/ul	95
12) Phenanthrene	17.29	178	278286	11.465	ng/ul#	91
13) Anthracene	17.38	178	51462	2.148	ng/ul#	91
15) Fluoranthene	19.31	202	480491	17.122	ng/ul	83
17) Pyrene	19.67	202	499661	17.674	ng/ul#	80
18) Benzo(a)anthracene	21.41	228	279254	10.682	ng/ul#	88
19) Chrysene	21.46	228	303643	11.237	ng/ul	99
21) Benzo(b)fluoranthene	23.18	252	297384m	11.157	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	101889m	3.567	ng/ul	
23) Benzo(a)pyrene	23.85	252	216820	8.137	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	26.67	276	111218	4.034	ng/ul#	75
25) Dibenzo(a,h)anthracene	26.66	278	29102	1.253	ng/ul	99
26) Benzo(g,h,i)perylene	27.52	276	83592	3.776	ng/ul	95

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

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