

Data Path : Z:\HPCHEM1\BNA E\DATA\BE112217\
 Data File : BE094843.D
 Acq On : 22 Nov 2017 21:53
 Operator : SJ/JU
 Sample : I6496-03
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AA2

Manual Integrations
 APPROVED

Sohil
 11/27/2017 3:25:43 PM

Quant Time: Nov 23 03:07:30 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE111417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 23 02:50:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	633	0.40	ng/ul	-0.04
2) Naphthalene-d8	10.68	136	4285	0.40	ng/ul	-0.03
6) Acenaphthene-d10	14.50	164	3086	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.24	188	7200	0.40	ng/ul	-0.02
16) Chrysene-d12	21.42	240	7706	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	8583	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.30	152	1775	0.26	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	6225	0.25	ng/ul	-0.02
Target Compounds					Ovalue	
3) Naphthalene	10.72	128	2223	0.218	ng/ul#	95
5) 2-Methylnaphthalene	12.37	142	274	0.039	ng/ul	98
7) Acenaphthylene	14.23	152	1396	0.099	ng/ul#	89
8) Acenaphthene	14.56	153	376	0.038	ng/ul#	88
9) Fluorene	15.55	166	682	0.057	ng/ul#	84
12) Phenanthrene	17.27	178	12422m	0.642	ng/ul	
13) Anthracene	17.38	178	3461	0.176	ng/ul	95
15) Fluoranthene	19.30	202	48359	1.906	ng/ul	84
17) Pyrene	19.66	202	45963	1.906	ng/ul#	79
18) Benzo(a)anthracene	21.40	228	28211	1.204	ng/ul#	89
19) Chrysene	21.45	228	25039	1.071	ng/ul	98
21) Benzo(b)fluoranthene	23.17	252	36692m	1.435	ng/ul	
22) Benzo(k)fluoranthene	23.21	252	11446m	0.421	ng/ul	
23) Benzo(a)pyrene	23.83	252	26548	1.044	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	26.65	276	17408	0.628	ng/ul#	72
25) Dibenzo(a,h)anthracene	26.64	278	4290	0.184	ng/ul	96
26) Benzo(a,h,i)perylene	27.49	276	14462	0.623	ng/ul	95

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

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