

Data Path : Z:\HPCHEM1\BNA E\DATA\BE112217\
 Data File : BE094854.D
 Acq On : 23 Nov 2017 4:21
 Operator : SJ/JU
 Sample : I6496-15
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AA6

Manual Integrations
 APPROVED

Sohil
 11/27/2017 3:26:11 PM

Quant Time: Nov 23 22:43:40 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	699	0.40	ng/ul	-0.04
2) Naphthalene-d8	10.67	136	4966	0.40	ng/ul	-0.03
6) Acenaphthene-d10	14.50	164	3539	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.24	188	8088	0.40	ng/ul	-0.02
16) Chrysene-d12	21.42	240	8242	0.40	ng/ul	-0.01
20) Perylene-d12	23.96	264	7422m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	2156	0.27	ng/ul	-0.01
14) Fluoranthene-d10	19.27	212	7428	0.27	ng/ul	-0.02
Target Compounds					Ovalue	
3) Naphthalene	10.72	128	383	0.032	ng/ul#	61
5) 2-Methylnaphthalene	12.37	142	228	0.028	ng/ul	92
7) Acenaphthylene	14.23	152	448	0.028	ng/ul#	66
8) Acenaphthene	14.56	153	362	0.032	ng/ul#	86
9) Fluorene	15.55	166	480	0.035	ng/ul#	87
12) Phenanthrene	17.27	178	10248	0.472	ng/ul#	90
13) Anthracene	17.36	178	2300	0.104	ng/ul#	92
15) Fluoranthene	19.30	202	21154	0.742	ng/ul	83
17) Pyrene	19.66	202	20530	0.796	ng/ul#	78
18) Benzo(a)anthracene	21.40	228	9811	0.392	ng/ul#	87
19) Chrysene	21.46	228	8826	0.353	ng/ul	98
21) Benzo(b)fluoranthene	23.17	252	10388m	0.470	ng/ul	
22) Benzo(k)fluoranthene	23.21	252	3776m	0.161	ng/ul	
23) Benzo(a)pyrene	23.83	252	8047m	0.366	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.66	276	4845m	0.202	ng/ul	
25) Dibenzo(a,h)anthracene	26.65	278	1218m	0.060	ng/ul	
26) Benzo(a,h,i)perylene	27.51	276	4723m	0.235	ng/ul	

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

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