

Data Path : Z:\HPCHEM1\BNA E\DATA\BE112217\
 Data File : BE094845.D
 Acq On : 22 Nov 2017 23:02
 Operator : SJ/JU
 Sample : I6496-05
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AB5

Manual Integrations
 APPROVED

Sohil
 11/27/2017 3:25:58 PM

Quant Time: Nov 23 03:12:20 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	780	0.40	ng/ul	-0.04
2) Naphthalene-d8	10.68	136	5396	0.40	ng/ul	-0.03
6) Acenaphthene-d10	14.50	164	3747	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.24	188	8829	0.40	ng/ul	-0.02
16) Chrysene-d12	21.41	240	8926	0.40	ng/ul	-0.01
20) Perylene-d12	23.95	264	9713	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	1716	0.20	ng/ul	-0.02
14) Fluoranthene-d10	19.27	212	5307	0.17	ng/ul	-0.02
Target Compounds				Ovalue		
3) Naphthalene	10.72	128	1287	0.100	ng/ul#	90
5) 2-Methylnaphthalene	12.36	142	889	0.099	ng/ul	99
7) Acenaphthylene	14.23	152	1821	0.106	ng/ul#	91
8) Acenaphthene	14.56	153	404	0.033	ng/ul#	91
9) Fluorene	15.55	166	735	0.050	ng/ul#	86
12) Phenanthrene	17.27	178	18846m	0.795	ng/ul	
13) Anthracene	17.38	178	3631	0.150	ng/ul	95
15) Fluoranthene	19.30	202	60890	1.957	ng/ul	82
17) Pyrene	19.66	202	52389	1.876	ng/ul#	78
18) Benzo(a)anthracene	21.40	228	31090	1.146	ng/ul#	88
19) Chrysene	21.45	228	35133	1.297	ng/ul	100
21) Benzo(b)fluoranthene	23.17	252	50301m	1.738	ng/ul	
22) Benzo(k)fluoranthene	23.21	252	16992m	0.552	ng/ul	
23) Benzo(a)pyrene	23.83	252	31119	1.081	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	26.64	276	23418	0.747	ng/ul#	72
25) Dibenzo(a,h)anthracene	26.64	278	5951	0.225	ng/ul#	93
26) Benzo(a,h,i)perylene	27.48	276	18249	0.694	ng/ul	96

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

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