

Data Path : Z:\HPCHEM1\BNA E\DATA\BE112217\
 Data File : BE094859.D
 Acq On : 23 Nov 2017 7:17
 Operator : SJ/JU
 Sample : I6496-19
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AB0

Manual Integrations
 APPROVED

Sohil
 11/27/2017 3:26:17 PM

Quant Time: Nov 23 23:48:33 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 23:11:00 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	689	0.40	ng/ul	-0.03
2) Naphthalene-d8	10.68	136	4796	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.50	164	3428	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.24	188	7926	0.40	ng/ul	-0.02
16) Chrysene-d12	21.42	240	8367	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	9245	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	1791	0.23	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	6110	0.22	ng/ul	-0.01
Target Compounds				Ovalue		
3) Naphthalene	10.72	128	298	0.026	ng/ul#	50
5) 2-Methylnaphthalene	12.37	142	212	0.027	ng/ul	97
7) Acenaphthylene	14.23	152	3113	0.198	ng/ul	99
12) Phenanthrene	17.27	178	5467	0.257	ng/ul#	89
13) Anthracene	17.36	178	2409	0.111	ng/ul#	92
15) Fluoranthene	19.30	202	44618	1.597	ng/ul	83
17) Pyrene	19.66	202	42870	1.637	ng/ul#	77
18) Benzo(a)anthracene	21.40	228	28034	1.102	ng/ul#	87
19) Chrysene	21.46	228	21832	0.860	ng/ul	99
21) Benzo(b)fluoranthene	23.17	252	35586m	1.292	ng/ul	
22) Benzo(k)fluoranthene	23.22	252	11439m	0.390	ng/ul	
23) Benzo(a)pyrene	23.83	252	21265	0.776	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	26.66	276	13447	0.450	ng/ul#	73
25) Dibenzo(a,h)anthracene	26.64	278	3514	0.140	ng/ul	97
26) Benzo(g,h,i)perylene	27.50	276	10343	0.413	ng/ul	95

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

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