

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110317\
 Data File : BE094716.D
 Acq On : 3 Nov 2017 20:19
 Operator : SJ/JU
 Sample : I5902-01 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-5(0.5-1.5)-101317

Manual Integrations
 APPROVED

Sohil
 11/4/2017 2:15:36 PM

Quant Time: Nov 04 00:25:00 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 04 01:15:30 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1472	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.69	136	8779	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	4819	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	10251	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	10285	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	11052	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	741	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	1694	0.06	ng/ul	0.00
Target Compounds						Qvalue
7) Acenaphthylene	14.24	152	458	0.021	ng/ul#	72
8) Acenaphthene	14.57	153	385	0.025	ng/ul#	94
9) Fluorene	15.56	166	409	0.023	ng/ul#	86
12) Phenanthrene	17.29	178	9566	0.353	ng/ul#	90
13) Anthracene	17.38	178	2547	0.095	ng/ul#	95
15) Fluoranthene	19.31	202	24012	0.767	ng/ul	84
17) Pyrene	19.67	202	22454	0.706	ng/ul#	82
18) Benzo(a)anthracene	21.40	228	11946	0.406	ng/ul#	88
19) Chrysene	21.46	228	12029	0.396	ng/ul	95
21) Benzo(b)fluoranthene	23.18	252	14960m	0.481	ng/ul	
22) Benzo(k)fluoranthene	23.22	252	6165m	0.185	ng/ul	
23) Benzo(a)pyrene	23.84	252	10293	0.331	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	26.66	276	6144	0.191	ng/ul#	74
25) Dibenzo(a,h)anthracene	26.65	278	1346	0.050	ng/ul#	21
26) Benzo(g,h,i)perylene	27.51	276	5287	0.205	ng/ul#	77

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

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