

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110317\
 Data File : BE094711.D
 Acq On : 3 Nov 2017 17:23
 Operator : SJ/JU
 Sample : I5825-09 5X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-64(1.5-2.5)-101117

Manual Integrations
 APPROVED

Sohil
 11/4/2017 2:15:22 PM

Quant Time: Nov 03 23:40:25 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 03 23:54:30 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1493	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.69	136	8815	0.40	ng/ul	-0.03
6) Acenaphthene-d10	14.51	164	4992	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.24	188	10679	0.40	ng/ul	-0.04
16) Chrysene-d12	21.42	240	10542	0.40	ng/ul	-0.02
20) Perylene-d12	23.95	264	11811	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	784	0.05	ng/ul	-0.02
14) Fluoranthene-d10	19.27	212	1862	0.06	ng/ul	-0.02
Target Compounds				Qvalue		
3) Naphthalene	10.74	128	4005	0.186	ng/ul#	96
5) 2-Methylnaphthalene	12.35	142	2103	0.139	ng/ul	97
7) Acenaphthylene	14.24	152	13461	0.596	ng/ul	97
8) Acenaphthene	14.57	153	1686	0.104	ng/ul#	94
9) Fluorene	15.56	166	4252	0.233	ng/ul	94
12) Phenanthrene	17.29	178	66941	2.373	ng/ul#	93
13) Anthracene	17.38	178	17677	0.635	ng/ul#	94
15) Fluoranthene	19.30	202	165860	5.086	ng/ul	84
17) Pyrene	19.66	202	150440	4.616	ng/ul#	81
18) Benzo(a)anthracene	21.40	228	91996	3.053	ng/ul#	87
19) Chrysene	21.45	228	90440	2.903	ng/ul	98
21) Benzo(b)fluoranthene	23.17	252	117150m	3.526	ng/ul	
22) Benzo(k)fluoranthene	23.22	252	43728m	1.228	ng/ul	
23) Benzo(a)pyrene	23.84	252	82047	2.471	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	26.65	276	51781	1.507	ng/ul#	74
25) Dibenzo(a,h)anthracene	26.64	278	12757	0.441	ng/ul#	82
26) Benzo(g,h,i)perylene	27.49	276	42227	1.530	ng/ul	98

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

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