

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110317\
 Data File : BE094705.D
 Acq On : 3 Nov 2017 13:48
 Operator : SJ/JU
 Sample : I5825-02 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-63(0-1)-101117

Manual Integrations
 APPROVED

Sohil
 11/4/2017 2:15:07 PM

Quant Time: Nov 03 23:17:38 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	1344	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.68	136	7862	0.40	ng/ul	-0.03
6) Acenaphthene-d10	14.51	164	4490	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.24	188	9359	0.40	ng/ul	-0.04
16) Chrysene-d12	21.42	240	9259	0.40	ng/ul	-0.02
20) Perylene-d12	23.95	264	10383	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	681	0.05	ng/ul	-0.03
14) Fluoranthene-d10	19.27	212	1386	0.05	ng/ul	-0.02
Target Compounds				Qvalue		
3) Naphthalene	10.73	128	17189	0.894	ng/ul#	94
5) 2-Methylnaphthalene	12.34	142	3513	0.261	ng/ul	95
7) Acenaphthylene	14.24	152	4334	0.213	ng/ul#	90
8) Acenaphthene	14.57	153	7768	0.532	ng/ul	98
9) Fluorene	15.56	166	9152	0.558	ng/ul	93
12) Phenanthrene	17.29	178	137906	5.579	ng/ul#	93
13) Anthracene	17.38	178	34815	1.427	ng/ul#	93
15) Fluoranthene	19.30	202	322803	11.295	ng/ul	84
17) Pyrene	19.66	202	281853	9.847	ng/ul#	80
18) Benzo(a)anthracene	21.40	228	152736	5.770	ng/ul#	87
19) Chrysene	21.46	228	146535	5.356	ng/ul	100
21) Benzo(b)fluoranthene	23.17	252	198572m	6.799	ng/ul	
22) Benzo(k)fluoranthene	23.21	252	69722m	2.228	ng/ul	
23) Benzo(a)pyrene	23.84	252	138665	4.750	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	26.65	276	81415	2.695	ng/ul#	75
25) Dibenzo(a,h)anthracene	26.64	278	17987	0.707	ng/ul	95
26) Benzo(g,h,i)perylene	27.48	276	66557	2.744	ng/ul	97

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

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