

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110317\
 Data File : BE094706.D
 Acq On : 3 Nov 2017 14:23
 Operator : SJ/JU
 Sample : I5825-03 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-63(1-3)-101117

Manual Integrations
 APPROVED

Sohil
 11/4/2017 2:15:10 PM

Quant Time: Nov 03 23:28:41 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	1503	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.68	136	8731	0.40	ng/ul	-0.03
6) Acenaphthene-d10	14.51	164	5036	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.24	188	10774	0.40	ng/ul	-0.04
16) Chrysene-d12	21.42	240	10288	0.40	ng/ul	-0.02
20) Perylene-d12	23.95	264	10363	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	561	0.04	ng/ul	-0.03
14) Fluoranthene-d10	19.27	212	1463	0.05	ng/ul	-0.02
Target Compounds				Qvalue		
3) Naphthalene	10.73	128	55577	2.603	ng/ul#	91
5) 2-Methylnaphthalene	12.34	142	21211	1.420	ng/ul	98
7) Acenaphthylene	14.24	152	4178	0.183	ng/ul#	93
8) Acenaphthene	14.57	153	152526	9.318	ng/ul	97
9) Fluorene	15.56	166	204996	11.153	ng/ul#	96
12) Phenanthrene	17.29	178	2289646	80.458	ng/ul#	90
13) Anthracene	17.38	178	636929	22.680	ng/ul#	91
15) Fluoranthene	19.31	202	2956885	89.874	ng/ul	83
17) Pyrene	19.67	202	2440439	76.732	ng/ul#	79
18) Benzo(a)anthracene	21.40	228	1172287	39.859	ng/ul#	87
19) Chrysene	21.46	228	1018179	33.493	ng/ul	99
21) Benzo(b)fluoranthene	23.18	252	1118952m	38.385	ng/ul	
22) Benzo(k)fluoranthene	23.22	252	397784m	12.734	ng/ul	
23) Benzo(a)pyrene	23.84	252	839983	28.827	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	26.64	276	410117	13.602	ng/ul#	76
25) Dibenzo(a,h)anthracene	26.63	278	94297	3.711	ng/ul	93
26) Benzo(g,h,i)perylene	27.48	276	331631	13.698	ng/ul	96

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

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