

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110317\
 Data File : BE094732.D
 Acq On : 4 Nov 2017 5:46
 Operator : SJ/JU
 Sample : I5933-01 5X
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-9(0-1) 101617

Manual Integrations
 APPROVED

Sohil
 11/4/2017 2:17:09 PM

Quant Time: Nov 04 05:45:01 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 06:43:16 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1706	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.68	136	10046	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.51	164	5606	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.24	188	12276	0.40	ng/ul	-0.02
16) Chrysene-d12	21.42	240	11966	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	11701	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	756	0.05	ng/ul	-0.02
14) Fluoranthene-d10	19.27	212	2082	0.06	ng/ul	-0.01
Target Compounds					Qvalue	
3) Naphthalene	10.73	128	53505	2.178	ng/ul#	91
5) 2-Methylnaphthalene	12.34	142	12904	0.751	ng/ul	97
7) Acenaphthylene	14.24	152	24983	0.985	ng/ul	97
8) Acenaphthene	14.57	153	56865	3.121	ng/ul	97
9) Fluorene	15.56	166	59858	2.925	ng/ul#	95
12) Phenanthrene	17.29	178	1085176	33.467	ng/ul#	92
13) Anthracene	17.38	178	208924	6.529	ng/ul#	91
15) Fluoranthene	19.31	202	1720737	45.902	ng/ul	83
17) Pyrene	19.67	202	1590167	42.987	ng/ul#	80
18) Benzo(a)anthracene	21.40	228	786009	22.978	ng/ul#	87
19) Chrysene	21.46	228	769677	21.768	ng/ul	99
21) Benzo(b)fluoranthene	23.18	252	881480m	26.781	ng/ul	
22) Benzo(k)fluoranthene	23.22	252	327091m	9.273	ng/ul	
23) Benzo(a)pyrene	23.85	252	639171	19.427	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	26.66	276	381156	11.196	ng/ul#	75
25) Dibenzo(a,h)anthracene	26.65	278	86493	3.015	ng/ul	94
26) Benzo(g,h,i)perylene	27.50	276	326661	11.949	ng/ul	95

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

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