

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110317\
 Data File : BE094722.D
 Acq On : 3 Nov 2017 23:54
 Operator : SJ/JU
 Sample : I5955-09 5X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-14(0-1)-101717

Manual Integrations
 APPROVED

Sohil
 11/4/2017 2:16:50 PM

Quant Time: Nov 04 00:41:26 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	1588	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.69	136	9229	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	5215	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.24	188	11288	0.40	ng/ul	-0.02
16) Chrysene-d12	21.42	240	10908	0.40	ng/ul	-0.01
20) Perylene-d12	23.95	264	11113	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	516	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	1315	0.04	ng/ul	-0.01
Target Compounds				Qvalue		
3) Naphthalene	10.74	128	5941	0.263	ng/ul#	94
5) 2-Methylnaphthalene	12.35	142	2057	0.130	ng/ul	97
7) Acenaphthylene	14.24	152	5076	0.215	ng/ul	98
8) Acenaphthene	14.57	153	8037	0.474	ng/ul	99
9) Fluorene	15.56	166	10386	0.546	ng/ul#	92
12) Phenanthrene	17.29	178	150049	5.033	ng/ul#	92
13) Anthracene	17.38	178	38085	1.294	ng/ul#	92
15) Fluoranthene	19.30	202	250874	7.278	ng/ul	84
17) Pyrene	19.66	202	218696	6.485	ng/ul#	81
18) Benzo(a)anthracene	21.40	228	117088	3.755	ng/ul#	87
19) Chrysene	21.45	228	118365	3.672	ng/ul	99
21) Benzo(b)fluoranthene	23.18	252	148911m	4.764	ng/ul	
22) Benzo(k)fluoranthene	23.22	252	46567m	1.390	ng/ul	
23) Benzo(a)pyrene	23.84	252	100715	3.223	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	26.65	276	56497	1.747	ng/ul#	74
25) Dibenzo(a,h)anthracene	26.65	278	14381	0.528	ng/ul#	78
26) Benzo(g,h,i)perylene	27.50	276	48033	1.850	ng/ul	97

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

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