

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110317\
 Data File : BE094740.D
 Acq On : 4 Nov 2017 10:29
 Operator : SJ/JU
 Sample : I5955-07MSD 5X
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-15(0.25-1.25)-101717MSD

Manual Integrations
 APPROVED

Sohil
 11/6/2017 5:58:52 PM

Quant Time: Nov 06 00:28:10 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 06:57:51 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	10.69	136	8938	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.51	164	5084	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	10838	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	11158	0.40	ng/ul	0.00
20) Perylene-d12	23.98	264	12036	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	1055	0.07	ng/ul	-0.02
14) Fluoranthene-d10	19.28	212	2475	0.08	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.74	128	7151	0.327	ng/ul#	94
5) 2-Methylnaphthalene	12.34	142	4413	0.289	ng/ul	98
7) Acenaphthylene	14.24	152	1960	0.085	ng/ul#	83
8) Acenaphthene	14.57	153	1529	0.093	ng/ul#	93
12) Phenanthrene	17.29	178	17401	0.608	ng/ul	92
13) Anthracene	17.38	178	2449m	0.087	ng/ul	
15) Fluoranthene	19.31	202	28721	0.868	ng/ul	82
17) Pyrene	19.67	202	41166	1.193	ng/ul#	84
18) Benzo(a)anthracene	21.41	228	16154	0.506	ng/ul#	42
19) Chrysene	21.47	228	37243	1.130	ng/ul#	60
21) Benzo(b)fluoranthene	23.19	252	29618m	0.875	ng/ul	
22) Benzo(k)fluoranthene	23.24	252	8360m	0.230	ng/ul	
23) Benzo(a)pyrene	23.86	252	16802	0.496	ng/ul#	45
24) Indeno(1,2,3-cd)pyrene	26.69	276	12727	0.363	ng/ul#	75
25) Dibenzo(a,h)anthracene	26.68	278	3620	0.123	ng/ul#	1
26) Benzo(g,h,i)perylene	27.55	276	14330	0.510	ng/ul#	59

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

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