

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE101317\
 Data File : BE094347.D
 Acq On : 14 Oct 2017 5:47
 Operator : SJ/JU
 Sample : I5649-03DL 5X
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-59(FILL)-100317DL

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:23:21 PM

Quant Time: Oct 14 22:29:35 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 14 05:22:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	2012	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	10259	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	5658	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	12082	0.40	ng/ul	-0.02
16) Chrysene-d12	21.43	240	12042	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	11452	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	392	0.02	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	969	0.02	ng/ul	0.00
Target Compounds						
						Qvalue
3) Naphthalene	10.76	128	2470	0.10	ng/ul#	93
5) 2-Methylnaphthalene	12.37	142	686	0.04	ng/ul	97
7) Acenaphthylene	14.25	152	2159	0.10	ng/ul	94
12) Phenanthrene	17.31	178	14219	0.45	ng/ul#	93
13) Anthracene	17.40	178	3507	0.11	ng/ul	95
15) Fluoranthene	19.32	202	40143	0.97	ng/ul	84
17) Pyrene	19.68	202	37708	1.13	ng/ul#	82
18) Benzo(a)anthracene	21.41	228	21775	0.68	ng/ul#	88
19) Chrysene	21.47	228	21912m	0.59	ng/ul	
21) Benzo(b)fluoranthene	23.19	252	30537m	0.98	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	10043m	0.28	ng/ul	
23) Benzo(a)pyrene	23.85	252	20822	0.64	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	26.65	276	15544	0.47	ng/ul#	78
25) Dibenzo(a,h)anthracene	26.65	278	4130	0.15	ng/ul#	24
26) Benzo(g,h,i)perylene	27.49	276	11601	0.41	ng/ul#	80

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

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