

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE101317\
 Data File : BE094350.D
 Acq On : 14 Oct 2017 7:38
 Operator : SJ/JU
 Sample : I5649-08DL 5X
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-57(1-2)-100317DL

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:23:27 PM

Quant Time: Oct 14 22:44:17 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	1947	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	10090	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	5568	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	12512	0.40	ng/ul	-0.02
16) Chrysene-d12	21.43	240	11743	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	12028	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	482	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	1227	0.03	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	3835	0.16	ng/ul#	95
5) 2-Methylnaphthalene	12.36	142	985	0.06	ng/ul	96
7) Acenaphthylene	14.26	152	3445	0.16	ng/ul	97
8) Acenaphthene	14.59	153	2529	0.14	ng/ul	97
9) Fluorene	15.57	166	3154	0.15	ng/ul	91
12) Phenanthrene	17.31	178	113303	3.47	ng/ul#	92
13) Anthracene	17.39	178	36405	1.14	ng/ul#	92
15) Fluoranthene	19.32	202	418221	9.73	ng/ul	84
17) Pyrene	19.68	202	388223	11.88	ng/ul#	80
18) Benzo(a)anthracene	21.41	228	257632	8.27	ng/ul#	87
19) Chrysene	21.47	228	219605	6.10	ng/ul	98
21) Benzo(b)fluoranthene	23.19	252	281205m	8.61	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	107315m	2.84	ng/ul	
23) Benzo(a)pyrene	23.85	252	224851	6.62	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	26.66	276	136349	3.93	ng/ul#	78
25) Dibenzo(a,h)anthracene	26.66	278	36195	1.24	ng/ul#	89
26) Benzo(g,h,i)perylene	27.50	276	106981	3.57	ng/ul	97

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

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