

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE062417\
 Data File : BE093359.D
 Acq On : 24 Jun 2017 13:25
 Operator : SJ/JU
 Sample : I3625-08
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5B82

Manual Integrations
 APPROVED

mohammad
 6/29/2017 8:34:07 AM

Quant Time: Jun 26 01:36:14 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 26 01:13:30 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	1417	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	7441	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	4883	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	11163m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	11827	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	9176m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.10	152	4175	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	13016	0.43	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	17796	0.964	ng/ul#	91
5) 2-Methylnaphthalene	12.18	142	4067	0.322	ng/ul	98
7) Acenaphthylene	14.08	152	79104	3.775	ng/ul	96
8) Acenaphthene	14.42	153	1540	0.091	ng/ul#	92
9) Fluorene	15.42	166	2493	0.126	ng/ul#	89
12) Phenanthrene	17.14	178	55002	1.791	ng/ul#	91
13) Anthracene	17.22	178	144112	5.389	ng/ul#	89
15) Fluoranthene	19.15	202	229080	6.082	ng/ul	83
17) Pyrene	19.51	202	248978	7.138	ng/ul#	85
18) Benzo(a)anthracene	21.24	228	196998	6.245	ng/ul#	82
19) Chrysene	21.29	228	225430	6.889	ng/ul	97
21) Benzo(b)fluoranthene	22.94	252	486297	17.463	ng/ul	88
22) Benzo(k)fluoranthene	22.98	252	146237m	4.945	ng/ul	
23) Benzo(a)pyrene	23.57	252	243756m	9.258	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.20	276	276575m	9.072	ng/ul	
25) Dibenzo(a,h)anthracene	26.20	278	65329	2.545	ng/ul	94
26) Benzo(g,h,i)perylene	26.97	276	270963	10.364	ng/ul	97

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

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