

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061917\
 Data File : BE093271.D
 Acq On : 20 Jun 2017 3:48
 Operator : SJ/MA
 Sample : I3599-06
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A87

Manual Integrations
 APPROVED

mohammad
 6/20/2017 3:13:31 PM

Quant Time: Jun 20 09:35:45 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 20 08:49:11 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	2528	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	13298	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	8440	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	19315	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	18794	0.40	ng/ul	0.01
20) Perylene-d12	23.68	264	24080m	0.40	ng/ul	0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	4569	0.23	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	11343	0.22	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	143593	4.352	ng/ul#	90
5) 2-Methylnaphthalene	12.18	142	30556	1.354	ng/ul	98
7) Acenaphthylene	14.09	152	197990	5.466	ng/ul	95
8) Acenaphthene	14.42	153	43092	1.475	ng/ul	98
9) Fluorene	15.42	166	43378	1.266	ng/ul	94
12) Phenanthrene	17.13	178	430044	8.092	ng/ul#	87
13) Anthracene	17.24	178	2967279	64.128	ng/ul#	94
15) Fluoranthene	19.17	202	6080101	93.292	ng/ul	84
17) Pyrene	19.54	202	6121191	110.442	ng/ul#	89
18) Benzo(a)anthracene	21.26	228	3429029	68.401	ng/ul#	84
19) Chrysene	21.31	228	3664957	70.476	ng/ul	95
21) Benzo(b)fluoranthene	22.96	252	4344391m	59.450	ng/ul	
22) Benzo(k)fluoranthene	23.00	252	1464447m	18.869	ng/ul	
23) Benzo(a)pyrene	23.59	252	2138486m	30.949	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.22	276	1286291m	16.077	ng/ul	
25) Dibenzo(a,h)anthracene	26.23	278	408099m	6.057	ng/ul	
26) Benzo(g,h,i)perylene	26.99	276	708564	10.327	ng/ul	96

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

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