

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE062317\
 Data File : BE093345.D
 Acq On : 23 Jun 2017 20:10
 Operator : SJ/JU
 Sample : I3625-11
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5D15

Manual Integrations
 APPROVED

mohammad
 6/29/2017 8:33:15 AM

Quant Time: Jun 24 04:48:55 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jun 24 04:29:07 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	1688	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	8973	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	5532	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	13565	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	13113	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	11001m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.10	152	2204	0.17	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	12248	0.33	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	3343	0.150	ng/ul#	96
5) 2-Methylnaphthalene	12.18	142	1626	0.107	ng/ul	100
7) Acenaphthylene	14.08	152	93297	3.930	ng/ul	96
8) Acenaphthene	14.42	153	5518	0.288	ng/ul	98
9) Fluorene	15.41	166	12953	0.577	ng/ul	92
12) Phenanthrene	17.14	178	439620	11.779	ng/ul#	90
13) Anthracene	17.22	178	134685	4.145	ng/ul#	91
15) Fluoranthene	19.15	202	1881599	41.109	ng/ul	84
17) Pyrene	19.52	202	1876769	48.532	ng/ul#	87
18) Benzo(a)anthracene	21.24	228	1124315	32.144	ng/ul#	84
19) Chrysene	21.30	228	1155390	31.843	ng/ul	98
21) Benzo(b)fluoranthene	22.94	252	1486772	44.534	ng/ul	88
22) Benzo(k)fluoranthene	22.98	252	593680m	16.743	ng/ul	
23) Benzo(a)pyrene	23.57	252	1023870m	32.435	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.20	276	705571m	19.304	ng/ul	
25) Dibenzo(a,h)anthracene	26.21	278	203813m	6.622	ng/ul	
26) Benzo(g,h,i)perylene	26.98	276	595388	18.994	ng/ul	97

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

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