

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE062317\
 Data File : BE093349.D
 Acq On : 23 Jun 2017 22:39
 Operator : SJ/JU
 Sample : I3627-09
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A69

Manual Integrations
 APPROVED

mohammad
 6/29/2017 8:33:31 AM

Quant Time: Jun 24 05:03:53 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	1655	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	8795	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	5586	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	13192	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	13887	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	10773m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	3108	0.24	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	8977	0.25	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	2334	0.107	ng/ul#	96
5) 2-Methylnaphthalene	12.18	142	691	0.046	ng/ul	100
7) Acenaphthylene	14.08	152	38742	1.616	ng/ul	96
8) Acenaphthene	14.42	153	909	0.047	ng/ul#	89
9) Fluorene	15.41	166	1849	0.082	ng/ul	88
12) Phenanthrene	17.13	178	31579	0.870	ng/ul#	91
13) Anthracene	17.22	178	44107	1.396	ng/ul#	90
15) Fluoranthene	19.15	202	334074	7.505	ng/ul	83
17) Pyrene	19.51	202	419878	10.253	ng/ul#	87
18) Benzo(a)anthracene	21.24	228	254943	6.882	ng/ul#	84
19) Chrysene	21.30	228	286395	7.453	ng/ul	98
21) Benzo(b)fluoranthene	22.93	252	470497	14.391	ng/ul	87
22) Benzo(k)fluoranthene	22.97	252	166464m	4.794	ng/ul	
23) Benzo(a)pyrene	23.56	252	244450m	7.908	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.19	276	164434m	4.594	ng/ul	
25) Dibenzo(a,h)anthracene	26.19	278	43298m	1.436	ng/ul	
26) Benzo(g,h,i)perylene	26.96	276	113437	3.696	ng/ul	97

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

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