

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE062317\
 Data File : BE093339.D
 Acq On : 23 Jun 2017 15:50
 Operator : SJ/JU
 Sample : I3625-06
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5B06

Manual Integrations
 APPROVED

mohammad
 6/29/2017 8:32:52 AM

Quant Time: Jun 24 04:27:34 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 03:58:33 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	1664	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	8737	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	5548	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	13252	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	13927	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	14536	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	2475	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	6936	0.19	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	1973	0.091	ng/ul#	95
5) 2-Methylnaphthalene	12.18	142	562	0.038	ng/ul	98
7) Acenaphthylene	14.09	152	17126	0.719	ng/ul	95
8) Acenaphthene	14.42	153	902	0.047	ng/ul	96
9) Fluorene	15.42	166	1041	0.046	ng/ul#	91
12) Phenanthrene	17.14	178	10073	0.276	ng/ul#	92
13) Anthracene	17.22	178	26127	0.823	ng/ul#	90
15) Fluoranthene	19.15	202	96123	2.150	ng/ul	84
17) Pyrene	19.51	202	150192	3.657	ng/ul#	84
18) Benzo(a)anthracene	21.24	228	71663	1.929	ng/ul#	85
19) Chrysene	21.29	228	98933	2.567	ng/ul	98
21) Benzo(b)fluoranthene	22.93	252	189030m	4.285	ng/ul	
22) Benzo(k)fluoranthene	22.97	252	68776m	1.468	ng/ul	
23) Benzo(a)pyrene	23.55	252	90757	2.176	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	26.18	276	62099	1.286	ng/ul#	87
25) Dibenzo(a,h)anthracene	26.19	278	16561	0.407	ng/ul#	85
26) Benzo(g,h,i)perylene	26.96	276	46070	1.112	ng/ul	96

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

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