

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062017\
 Data File : BE093298.D
 Acq On : 21 Jun 2017 1:20
 Operator : SJ/JU
 Sample : I3627-11
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A99

Manual Integrations
 APPROVED

mohammad
 6/21/2017 12:19:11 PM

Quant Time: Jun 21 04:10:15 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 21 03:23:49 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	1717	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	8955	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	5706	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	13645	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	13501	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	10516m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	3526	0.27	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	10367	0.28	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	2535	0.114	ng/ul#	96
5) 2-Methylnaphthalene	12.18	142	1306	0.086	ng/ul	99
7) Acenaphthylene	14.09	152	6201	0.253	ng/ul	98
8) Acenaphthene	14.42	153	708	0.036	ng/ul	96
9) Fluorene	15.42	166	4883	0.211	ng/ul	95
12) Phenanthrene	17.14	178	17760	0.473	ng/ul#	89
13) Anthracene	17.22	178	99648	3.048	ng/ul#	89
15) Fluoranthene	19.15	202	48261	1.048	ng/ul	84
17) Pyrene	19.51	202	50280	1.263	ng/ul#	87
18) Benzo(a)anthracene	21.24	228	26380	0.733	ng/ul#	85
19) Chrysene	21.29	228	30348	0.812	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	38179	1.196	ng/ul	89
22) Benzo(k)fluoranthene	22.97	252	13653m	0.403	ng/ul	
23) Benzo(a)pyrene	23.55	252	20108m	0.666	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.18	276	13103m	0.375	ng/ul	
25) Dibenzo(a,h)anthracene	26.19	278	3460m	0.118	ng/ul	
26) Benzo(g,h,i)perylene	26.95	276	9235	0.308	ng/ul#	94

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

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