

Data Path : Z:\HPCHEM1\BNA E\DATA\BE062117\
 Data File : BE093313.D
 Acq On : 21 Jun 2017 13:40
 Operator : SJ/JU
 Sample : I3627-18DL 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A91DL

Manual Integrations
 APPROVED

mohammad
 6/22/2017 11:48:42 AM

Quant Time: Jun 22 03:16:40 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 22 02:10:17 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	2166	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	12101	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	7635	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	18685	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	20280	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	17648	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	628	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	1763	0.03	ng/ul	0.00
Target Compounds						Ovalue
3) Naphthalene	10.56	128	1880	0.06	ng/ul#	93
5) 2-Methylnaphthalene	12.18	142	1154	0.06	ng/ul	99
7) Acenaphthylene	14.09	152	4326	0.13	ng/ul	99
8) Acenaphthene	14.42	153	3277	0.12	ng/ul	98
9) Fluorene	15.41	166	4675	0.15	ng/ul	92
12) Phenanthrene	17.14	178	129444	2.52	ng/ul#	91
13) Anthracene	17.22	178	24365	0.54	ng/ul#	92
15) Fluoranthene	19.15	202	232476	3.69	ng/ul	84
17) Pyrene	19.51	202	220537	3.69	ng/ul#	83
18) Benzo(a)anthracene	21.24	228	107384	1.99	ng/ul#	85
19) Chrysene	21.29	228	110212	1.96	ng/ul	99
21) Benzo(b)fluoranthene	22.93	252	114250m	2.13	ng/ul	
22) Benzo(k)fluoranthene	22.97	252	40469m	0.71	ng/ul	
23) Benzo(a)pyrene	23.56	252	75269	1.49	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	26.18	276	43994	0.75	ng/ul#	87
25) Dibenzo(a,h)anthracene	26.19	278	13463	0.27	ng/ul	92
26) Benzo(a,h,i)perylene	26.95	276	37193	0.74	ng/ul	97

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

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