

Data Path : Z:\HPCHEM1\BNA E\DATA\BE062117\
 Data File : BE093322.D
 Acq On : 21 Jun 2017 19:11
 Operator : SJ/JU
 Sample : I3627-02DL 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A08DL

Manual Integrations
 APPROVED

mohammad
 6/22/2017 11:49:15 AM

Quant Time: Jun 22 03:46:02 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	7.74	152	1676	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	9304	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	5701	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	14151	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	15468	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	14598	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	819	0.06	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	2166	0.06	ng/ul	0.00
Target Compounds						Ovalue
3) Naphthalene	10.56	128	1001	0.04	ng/ul#	88
5) 2-Methylnaphthalene	12.18	142	248	0.02	ng/ul	96
7) Acenaphthylene	14.08	152	2973	0.12	ng/ul	98
8) Acenaphthene	14.42	153	750	0.04	ng/ul	97
9) Fluorene	15.41	166	1254	0.05	ng/ul	93
12) Phenanthrene	17.14	178	12879	0.33	ng/ul#	91
13) Anthracene	17.22	178	4348	0.13	ng/ul#	93
15) Fluoranthene	19.15	202	32140	0.67	ng/ul	84
17) Pyrene	19.51	202	40874	0.90	ng/ul#	88
18) Benzo(a)anthracene	21.24	228	20924	0.51	ng/ul#	86
19) Chrysene	21.29	228	26167	0.61	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	36485m	0.82	ng/ul	
22) Benzo(k)fluoranthene	22.97	252	14381m	0.31	ng/ul	
23) Benzo(a)pyrene	23.55	252	20530	0.49	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	26.18	276	12515	0.26	ng/ul#	88
25) Dibenzo(a,h)anthracene	26.19	278	3347	0.08	ng/ul#	81
26) Benzo(a,h,i)perylene	26.95	276	9012	0.22	ng/ul	96

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

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