

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061417\
 Data File : BE093184.D
 Acq On : 15 Jun 2017 00:05
 Operator : SJ/MA
 Sample : I3521-06
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5C54

Manual Integrations
 APPROVED

mohammad
 6/15/2017 3:53:40 PM

Quant Time: Jun 15 04:31:27 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 15 03:59:13 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	3991	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	20842	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	13049	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	29605	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	29462	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	23687m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	7436	0.24	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	19853	0.25	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.57	128	12815	0.248	ng/ul#	93
5) 2-Methylnaphthalene	12.18	142	3142	0.089	ng/ul	98
7) Acenaphthylene	14.09	152	40527	0.724	ng/ul	96
8) Acenaphthene	14.43	153	2244	0.050	ng/ul	96
9) Fluorene	15.42	166	1806	0.034	ng/ul#	82
12) Phenanthrene	17.14	178	390469	4.794	ng/ul#	86
13) Anthracene	17.22	178	73675	1.039	ng/ul#	87
15) Fluoranthene	19.16	202	697730	6.985	ng/ul	83
17) Pyrene	19.52	202	504758	5.810	ng/ul#	85
18) Benzo(a)anthracene	21.24	228	132538	1.687	ng/ul#	85
19) Chrysene	21.29	228	284462	3.489	ng/ul	97
21) Benzo(b)fluoranthene	22.94	252	314726	4.378	ng/ul	89
22) Benzo(k)fluoranthene	22.98	252	118821m	1.556	ng/ul	
23) Benzo(a)pyrene	23.56	252	103864m	1.528	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.19	276	104977m	1.334	ng/ul	
25) Dibenzo(a,h)anthracene	26.20	278	23754m	0.358	ng/ul	
26) Benzo(g,h,i)perylene	26.96	276	84258	1.248	ng/ul	95

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

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