

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061317\
 Data File : BE093152.D
 Acq On : 14 Jun 2017 4:18
 Operator : SJ/MA
 Sample : I3520-11
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A43

Manual Integrations
 APPROVED

mohammad
 6/14/2017 2:59:59 PM

Quant Time: Jun 14 11:22:47 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 14 08:39:04 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	4846	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	24566	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	14818	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	35555	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	37922	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	30148m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	8015	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	22180	0.23	ng/ul	0.00
Target Compounds				Qvalue		
12) Phenanthrene	17.14	178	6131	0.063	ng/ul#	90
15) Fluoranthene	19.15	202	17216	0.144	ng/ul	84
17) Pyrene	19.51	202	19597	0.175	ng/ul#	91
18) Benzo(a)anthracene	21.25	228	8642	0.085	ng/ul#	89
19) Chrysene	21.30	228	10313m	0.098	ng/ul	
21) Benzo(b)fluoranthene	22.93	252	15369	0.168	ng/ul	87
22) Benzo(k)fluoranthene	22.97	252	6000m	0.062	ng/ul	
23) Benzo(a)pyrene	23.56	252	8909m	0.103	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.18	276	6165m	0.062	ng/ul	
26) Benzo(g,h,i)perylene	26.95	276	5641	0.066	ng/ul#	83

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

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