

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061517\
 Data File : BE093210.D
 Acq On : 15 Jun 2017 16:12
 Operator : SJ/MA
 Sample : I3558-07DL 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5C61DL

Manual Integrations
 APPROVED

mohammad
 6/19/2017 11:06:22 AM

Quant Time: Jun 16 04:53:51 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 16 04:43:11 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	2928	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	16110	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	9767	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	25376	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	24924	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	34541	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	883	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	2443	0.04	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.57	128	14906	0.373	ng/ul#	92
5) 2-Methylnaphthalene	12.19	142	2954	0.108	ng/ul	99
7) Acenaphthylene	14.09	152	3150	0.075	ng/ul	98
8) Acenaphthene	14.43	153	4303	0.127	ng/ul	99
9) Fluorene	15.42	166	2824	0.071	ng/ul	93
12) Phenanthrene	17.13	178	22540	0.323	ng/ul#	86
13) Anthracene	17.22	178	8868	0.146	ng/ul#	88
15) Fluoranthene	19.16	202	34164	0.399	ng/ul	83
17) Pyrene	19.52	202	30228	0.411	ng/ul#	84
18) Benzo(a)anthracene	21.25	228	14964	0.225	ng/ul#	85
19) Chrysene	21.30	228	22400	0.325	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	59585m	0.568	ng/ul	
22) Benzo(k)fluoranthene	22.98	252	15307m	0.137	ng/ul	
23) Benzo(a)pyrene	23.56	252	28084	0.283	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	26.19	276	31250	0.272	ng/ul#	87
25) Dibenzo(a,h)anthracene	26.20	278	7461	0.077	ng/ul#	86
26) Benzo(g,h,i)perylene	26.96	276	28172	0.286	ng/ul	97

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

