

Data Path : Z:\HPCHEM1\BNA_E\Data\BE061317\
 Data File : BE093158.D
 Acq On : 14 Jun 2017 7:58
 Operator : SJ/MA
 Sample : I3520-03
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Jun 14 11:47:01 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	6094	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	30974	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	18319	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	43033	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	43533	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	45136	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	7886	0.17	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	19486	0.17	ng/ul	0.00
Target Compounds				Qvalue		
11) Pentachlorophenol	16.75	266	96	0.015	ng/ul	96
12) Phenanthrene	17.13	178	4698	0.040	ng/ul#	89
13) Anthracene	17.22	178	1187	0.012	ng/ul#	52
15) Fluoranthene	19.15	202	11131	0.077	ng/ul	83
17) Pyrene	19.51	202	13473	0.105	ng/ul#	92
18) Benzo(a)anthracene	21.24	228	4744	0.041	ng/ul#	86
19) Chrysene	21.29	228	5849	0.049	ng/ul	94
21) Benzo(b)fluoranthene	22.93	252	11960	0.087	ng/ul#	71
22) Benzo(k)fluoranthene	22.93	252	11960	0.082	ng/ul#	68
23) Benzo(a)pyrene	23.55	252	4631	0.036	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	26.18	276	3467	0.023	ng/ul#	60
26) Benzo(g,h,i)perylene	26.95	276	3043	0.024	ng/ul#	72

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

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