

Data Path : Z:\HPCHEM1\BNA_E\Data\BE061417\
 Data File : BE093200.D
 Acq On : 15 Jun 2017 9:53
 Operator : SJ/MA
 Sample : I3520-15
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5B47

Quant Time: Jun 15 11:33:04 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 07:58:47 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	3120	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	16435	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	10225	0.40	ng/ul	0.00
10) Phenanthrene-d10	0.00	188	0	0.00	ng/ul	-17.10
16) Chrysene-d12	21.27	240	23887	0.40	ng/ul	0.00
20) Perylene-d12	23.53	264	1143364	0.40	ng/ul	-0.13

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.11	152	7167	0.30	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	16829	0.00	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.57	128	18245	0.447	ng/ul#	92
5) 2-Methylnaphthalene	12.19	142	4367	0.157	ng/ul	98
7) Acenaphthylene	14.08	152	199508	4.547	ng/ul	95
8) Acenaphthene	14.43	153	5595	0.158	ng/ul	96
9) Fluorene	15.42	166	8005	0.193	ng/ul	90
17) Pyrene	19.53	202	4075052	57.848	ng/ul	92
18) Benzo(a)anthracene	21.25	228	1934077	30.355	ng/ul#	83
19) Chrysene	21.31	228	2736335	41.400	ng/ul	97
21) Benzo(b)fluoranthene	22.96	252	3343065	0.963	ng/ul	89
22) Benzo(k)fluoranthene	22.96	252	4726913	1.283	ng/ul#	91
23) Benzo(a)pyrene	23.47	252	1597936	0.487	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	25.93	276	167449	0.044	ng/ul#	75
25) Dibenzo(a,h)anthracene	26.11	278	32274	0.010	ng/ul	94
26) Benzo(g,h,i)perylene	26.99	276	747924	0.230	ng/ul	97

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

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