

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE112817\
 Data File : BE094881.D
 Acq On : 28 Nov 2017 13:24
 Operator : SJ/JU
 Sample : I6514-03DL 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AC3DL

Quant Time: Nov 29 01:13:28 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE111417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 29 01:10:42 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	621	0.40	ng/ul	-0.06
2) Naphthalene-d8	10.69	136	4333	0.40	ng/ul	-0.03
6) Acenaphthene-d10	14.50	164	2788	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.24	188	6777	0.40	ng/ul	-0.02
16) Chrysene-d12	21.43	240	6653	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	6924	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	454	0.06	ng/ul	-0.02
14) Fluoranthene-d10	19.28	212	1384	0.06	ng/ul	-0.01

Target Compounds				Qvalue		
3) Naphthalene	10.73	128	900	0.087	ng/ul#	83
5) 2-Methylnaphthalene	12.36	142	287	0.040	ng/ul	96
8) Acenaphthene	14.57	153	830	0.092	ng/ul	97
9) Fluorene	15.56	166	1196	0.110	ng/ul	93
12) Phenanthrene	17.27	178	7817	0.429	ng/ul#	89
13) Anthracene	17.38	178	1137	0.061	ng/ul#	88
15) Fluoranthene	19.30	202	2746	0.115	ng/ul	85
17) Pyrene	19.66	202	2748	0.132	ng/ul#	87
18) Benzo(a)anthracene	21.42	228	649	0.032	ng/ul#	76
19) Chrysene	21.47	228	659	0.033	ng/ul#	79
21) Benzo(b)fluoranthene	23.18	252	550	0.027	ng/ul#	63
23) Benzo(a)pyrene	23.84	252	436	0.021	ng/ul#	46
26) Benzo(g,h,i)perylene	27.59	276	378	0.020	ng/ul#	38

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

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