

Data Path : Z:\HPCHEM1\BNA E\DATA\BE112217\
 Data File : BE094833.D
 Acq On : 22 Nov 2017 15:58
 Operator : SJ/JU
 Sample : I6514-07
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AC8

Manual Integrations
 APPROVED

Sohil
 11/27/2017 3:25:07 PM

Quant Time: Nov 23 01:38:36 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE111417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 23 00:56:52 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	706	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.67	136	4979	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.50	164	3336	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.24	188	7675m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.42	240	7911	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	6763m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	2936	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	10078	0.38	ng/ul	-0.01
Target Compounds						
3) Naphthalene	10.72	128	449	0.038	ng/ul#	66
9) Fluorene	15.55	166	557	0.043	ng/ul#	90
12) Phenanthrene	17.27	178	3372	0.164	ng/ul	92
13) Anthracene	17.36	178	1255m	0.060	ng/ul	
15) Fluoranthene	19.30	202	2566	0.095	ng/ul#	52
17) Pyrene	19.66	202	1745	0.070	ng/ul#	92
18) Benzo(a)anthracene	21.40	228	894	0.037	ng/ul#	73
19) Chrysene	21.46	228	883	0.037	ng/ul#	74
21) Benzo(b)fluoranthene	23.18	252	761m	0.038	ng/ul	

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

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