

Data Path : Z:\HPCHEM1\BNA E\DATA\BE112217\
 Data File : BE094852.D
 Acq On : 23 Nov 2017 3:10
 Operator : SJ/JU
 Sample : I6496-13
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AA3

Manual Integrations
 APPROVED

Sohil
 11/27/2017 3:26:07 PM

Quant Time: Nov 23 03:53:43 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 02:50:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	715	0.40	ng/ul	-0.05
2) Naphthalene-d8	10.67	136	4821	0.40	ng/ul	-0.04
6) Acenaphthene-d10	14.50	164	3259	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.24	188	7930m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.42	240	8332	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	7522m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	2221	0.29	ng/ul	-0.03
14) Fluoranthene-d10	19.27	212	7927	0.29	ng/ul	-0.01
Target Compounds						
3) Naphthalene	10.72	128	1017	0.089	ng/ul#	84
11) Pentachlorophenol	16.96	266	127m	0.304	ng/ul	
12) Phenanthrene	17.27	178	1234m	0.058	ng/ul	
15) Fluoranthene	19.30	202	1571	0.056	ng/ul#	77
17) Pyrene	19.66	202	1350	0.052	ng/ul#	88
18) Benzo(a)anthracene	21.40	228	859	0.034	ng/ul#	72
19) Chrysene	21.45	228	1019	0.040	ng/ul#	70
21) Benzo(b)fluoranthene	23.18	252	1010m	0.045	ng/ul	
23) Benzo(a)pyrene	23.83	252	686m	0.031	ng/ul	

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

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