

Data Path : Z:\HPCHEM1\BNA E\DATA\BE101417\
 Data File : BE094384.D
 Acq On : 15 Oct 2017 8:54
 Operator : SJ/JU
 Sample : I5548-07DL 5X
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AB8DL

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:25:55 PM

Quant Time: Oct 16 06:46:55 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 05:09:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1663	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	8936	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	5426	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	12623	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	12621	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	12468	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	884	0.06	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	2332	0.06	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.29	178	6372	0.193	ng/ul#	92
13) Anthracene	17.38	178	1206m	0.038	ng/ul	
15) Fluoranthene	19.31	202	13911	0.321	ng/ul	83
17) Pyrene	19.67	202	12057	0.343	ng/ul#	83
18) Benzo(a)anthracene	21.40	228	6520	0.195	ng/ul#	68
21) Benzo(b)fluoranthene	23.17	252	9173m	0.271	ng/ul	
22) Benzo(k)fluoranthene	23.21	252	4237m	0.108	ng/ul	
23) Benzo(a)pyrene	23.83	252	6101	0.173	ng/ul#	64
24) Indeno(1,2,3-cd)pyrene	26.63	276	4152	0.115	ng/ul#	78

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

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