

Data Path : Z:\HPCHEM1\BNA E\DATA\BE081617\
 Data File : BE093856.D
 Acq On : 17 Aug 2017 1:01
 Operator : SJ/JU
 Sample : I4672-10
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AJ0

Manual Integrations
 APPROVED

Sohil
 8/17/2017 5:59:53 PM

Quant Time: Aug 17 04:56:06 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 17 02:52:40 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2475	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	13021	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	7737	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	15937	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	17211	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	15945m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	4679	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	12590	0.25	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.75	128	6582	0.208	ng/ul#	94
5) 2-Methylnaphthalene	12.36	142	4806	0.208	ng/ul	98
7) Acenaphthylene	14.25	152	6146	0.190	ng/ul	98
8) Acenaphthene	14.59	153	7529	0.296	ng/ul	99
9) Fluorene	15.57	166	7351	0.228	ng/ul	95
12) Phenanthrene	17.29	178	138069	3.181	ng/ul#	87
13) Anthracene	17.38	178	38637	0.914	ng/ul#	88
15) Fluoranthene	19.30	202	300452	5.535	ng/ul	84
17) Pyrene	19.67	202	271148	5.397	ng/ul#	81
18) Benzo(a)anthracene	21.39	228	160139	3.209	ng/ul#	86
19) Chrysene	21.45	228	163509	3.416	ng/ul	98
21) Benzo(b)fluoranthene	23.15	252	219247	4.445	ng/ul	87
22) Benzo(k)fluoranthene	23.19	252	87109m	1.791	ng/ul	
23) Benzo(a)pyrene	23.81	252	165498m	3.527	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.57	276	110506m	2.133	ng/ul	
25) Dibenzo(a,h)anthracene	26.57	278	26170m	0.605	ng/ul	
26) Benzo(a,h,i)perylene	27.39	276	104057m	2.384	ng/ul	

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

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