

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081517\
 Data File : BE093825.D
 Acq On : 16 Aug 2017 2:00
 Operator : SJ/JU
 Sample : I4672-20
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AG9

Manual Integrations
 APPROVED

Sohil
 8/16/2017 6:42:54 PM

Quant Time: Aug 16 08:24:45 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 16 07:30:49 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2023	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	10647	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	6677	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	14620	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	15615	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	16129m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	3677	0.21	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	10534	0.23	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	25150	0.971	ng/ul#	92
5) 2-Methylnaphthalene	12.36	142	22854	1.210	ng/ul	99
7) Acenaphthylene	14.24	152	4038	0.145	ng/ul#	87
9) Fluorene	15.57	166	999	0.036	ng/ul#	76
12) Phenanthrene	17.29	178	46692	1.173	ng/ul#	90
13) Anthracene	17.38	178	6687	0.172	ng/ul	93
15) Fluoranthene	19.30	202	72903	1.464	ng/ul	84
17) Pyrene	19.66	202	66048	1.449	ng/ul#	82
18) Benzo(a)anthracene	21.39	228	33969	0.750	ng/ul#	86
19) Chrysene	21.45	228	38401	0.884	ng/ul	97
21) Benzo(b)fluoranthene	23.14	252	44712	0.896	ng/ul	89
22) Benzo(k)fluoranthene	23.17	252	20002m	0.407	ng/ul	
23) Benzo(a)pyrene	23.80	252	29980m	0.632	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.55	276	18777m	0.358	ng/ul	
25) Dibenzo(a,h)anthracene	26.56	278	4449m	0.102	ng/ul	
26) Benzo(g,h,i)perylene	27.37	276	14536m	0.329	ng/ul	

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

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