

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081017\
 Data File : BE093749.D
 Acq On : 10 Aug 2017 15:52
 Operator : SJ/JU
 Sample : I4455-09
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AD5

Manual Integrations
 APPROVED

sohil
 8/14/2017 7:02:35 PM

Quant Time: Aug 11 18:08:55 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 11 06:07:04 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2923	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	14847	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	8933	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	20388	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	23190	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	19478m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	4973	0.21	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	14642	0.23	ng/ul	0.00
Target Compounds				Qvalue		
5) 2-Methylnaphthalene	12.36	142	1297	0.049	ng/ul	99
9) Fluorene	15.57	166	1435	0.039	ng/ul#	89
12) Phenanthrene	17.29	178	6320	0.114	ng/ul	92
15) Fluoranthene	19.30	202	21313	0.307	ng/ul	84
17) Pyrene	19.66	202	26394	0.390	ng/ul#	82
18) Benzo(a)anthracene	21.39	228	15034	0.224	ng/ul#	88
19) Chrysene	21.45	228	14868	0.231	ng/ul	98
21) Benzo(b)fluoranthene	23.14	252	17057m	0.283	ng/ul	
22) Benzo(k)fluoranthene	23.19	252	5824m	0.098	ng/ul	
23) Benzo(a)pyrene	23.80	252	13127m	0.229	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.54	276	7618m	0.120	ng/ul	
25) Dibenzo(a,h)anthracene	26.55	278	2127m	0.040	ng/ul	
26) Benzo(g,h,i)perylene	27.35	276	7530m	0.141	ng/ul	

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

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