

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072417\
 Data File : BE093560.D
 Acq On : 25 Jul 2017 4:30
 Operator : SJ/JU
 Sample : I4213-08
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0B32

Manual Integrations
 APPROVED

Sohil
 7/25/2017 5:33:58 PM

Quant Time: Jul 25 07:29:58 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 25 06:10:24 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	3290	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	17893	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	12043	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	28971m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	31523	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	28198m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	9400	0.32	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	30388	0.33	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.76	128	2133	0.049	ng/ul#	94
7) Acenaphthylene	14.25	152	1816	0.036	ng/ul#	90
12) Phenanthrene	17.29	178	4698m	0.060	ng/ul	
15) Fluoranthene	19.30	202	20761	0.210	ng/ul	84
17) Pyrene	19.67	202	22607	0.246	ng/ul#	82
18) Benzo(a)anthracene	21.39	228	12707	0.139	ng/ul#	86
19) Chrysene	21.45	228	12331	0.141	ng/ul	97
21) Benzo(b)fluoranthene	23.15	252	18841	0.216	ng/ul	93
22) Benzo(k)fluoranthene	23.20	252	6682m	0.078	ng/ul	
23) Benzo(a)pyrene	23.81	252	14122m	0.170	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.55	276	8644m	0.094	ng/ul	
26) Benzo(g,h,i)perylene	27.37	276	7891m	0.102	ng/ul	

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

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