

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072417\
 Data File : BE093569.D
 Acq On : 25 Jul 2017 10:04
 Operator : SJ/JU
 Sample : I4209-08
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0B11

Manual Integrations
 APPROVED

Sohil
 7/25/2017 5:34:15 PM

Quant Time: Jul 25 12:19:54 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	7.92	152	2219	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	11997	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	8151	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	20370m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	22336	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	18584m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	5568	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	18440	0.29	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.76	128	1893	0.065	ng/ul#	93
12) Phenanthrene	17.29	178	1697	0.031	ng/ul#	90
15) Fluoranthene	19.31	202	4137	0.060	ng/ul	85
17) Pyrene	19.67	202	4797m	0.074	ng/ul	
18) Benzo(a)anthracene	21.39	228	2698	0.042	ng/ul#	89
19) Chrysene	21.45	228	2985	0.048	ng/ul	93
21) Benzo(b)fluoranthene	23.15	252	4742	0.082	ng/ul	87
22) Benzo(k)fluoranthene	23.19	252	2042m	0.036	ng/ul	
23) Benzo(a)pyrene	23.80	252	3969m	0.073	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.55	276	2552m	0.042	ng/ul	
26) Benzo(g,h,i)perylene	27.37	276	2501m	0.049	ng/ul	

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

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