

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072417\
 Data File : BE093554.D
 Acq On : 25 Jul 2017 00:49
 Operator : SJ/JU
 Sample : I4213-12
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0B36

Manual Integrations
 APPROVED

Sohil
 7/25/2017 5:33:21 PM

Quant Time: Jul 25 11:21:47 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	2991	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	16100	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	10896	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	26801m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	29399	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	24322m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	8477	0.32	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	27837	0.33	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.76	128	13174	0.336	ng/ul#	92
5) 2-Methylnaphthalene	12.37	142	1282	0.045	ng/ul	99
12) Phenanthrene	17.29	178	2205	0.030	ng/ul#	90
13) Anthracene	17.40	178	2325	0.033	ng/ul#	86
15) Fluoranthene	19.31	202	7000	0.077	ng/ul	86
17) Pyrene	19.67	202	7671m	0.089	ng/ul	
18) Benzo(a)anthracene	21.39	228	5594	0.066	ng/ul#	86
19) Chrysene	21.45	228	6460	0.079	ng/ul	94
21) Benzo(b)fluoranthene	23.15	252	8982	0.119	ng/ul	92
22) Benzo(k)fluoranthene	23.19	252	4002m	0.054	ng/ul	
23) Benzo(a)pyrene	23.81	252	7792m	0.109	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.55	276	4664m	0.059	ng/ul	
26) Benzo(g,h,i)perylene	27.37	276	4055m	0.061	ng/ul	

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

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