

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072517\
 Data File : BE093585.D
 Acq On : 25 Jul 2017 19:55
 Operator : SJ/JU
 Sample : I4213-17
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0B41

Manual Integrations
 APPROVED

Sohil
 7/26/2017 5:56:00 PM

Quant Time: Jul 26 04:17:21 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	2372	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	12526	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	8460	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	20793m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	22739	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	19428m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	6743	0.33	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	22609	0.34	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.76	128	2630	0.086	ng/ul#	95
12) Phenanthrene	17.29	178	1705	0.030	ng/ul#	91
15) Fluoranthene	19.30	202	3501	0.049	ng/ul	86
17) Pyrene	19.66	202	3946	0.059	ng/ul#	92
18) Benzo(a)anthracene	21.39	228	2759	0.042	ng/ul#	88
19) Chrysene	21.45	228	2823m	0.045	ng/ul	
21) Benzo(b)fluoranthene	23.14	252	5287	0.088	ng/ul	90
22) Benzo(k)fluoranthene	23.19	252	2058m	0.035	ng/ul	
23) Benzo(a)pyrene	23.80	252	4118m	0.072	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.54	276	2768m	0.044	ng/ul	
26) Benzo(g,h,i)perylene	27.36	276	2634m	0.050	ng/ul	

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

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