

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102717\
 Data File : BE094598.D
 Acq On : 27 Oct 2017 16:26
 Operator : SJ/JU
 Sample : I5719-07 5X
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-25(0.5-1.5)-100517

Manual Integrations
 APPROVED

Sohil
 10/30/2017 6:36:51 PM

Quant Time: Oct 30 01:21:19 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 30 01:11:00 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1128	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	6409	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	3678	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	8141	0.40	ng/ul	-0.02
16) Chrysene-d12	21.43	240	7910	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	8961	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	498	0.05	ng/ul	-0.01
14) Fluoranthene-d10	19.29	212	1076	0.04	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.74	128	4597	0.293	ng/ul#	96
5) 2-Methylnaphthalene	12.35	142	2312	0.211	ng/ul	97
7) Acenaphthylene	14.24	152	5214	0.313	ng/ul#	93
8) Acenaphthene	14.58	153	10213	0.854	ng/ul	97
9) Fluorene	15.57	166	10778	0.803	ng/ul	92
12) Phenanthrene	17.29	178	341535	15.883	ng/ul#	87
13) Anthracene	17.39	178	86585	4.080	ng/ul	92
15) Fluoranthene	19.32	202	935177	37.618	ng/ul	84
17) Pyrene	19.68	202	860431	35.187	ng/ul#	81
18) Benzo(a)anthracene	21.41	228	656457	29.031	ng/ul#	87
19) Chrysene	21.47	228	598555	25.609	ng/ul	99
21) Benzo(b)fluoranthene	23.19	252	901610m	35.769	ng/ul	
22) Benzo(k)fluoranthene	23.24	252	286684m	10.613	ng/ul	
23) Benzo(a)pyrene	23.86	252	676677	26.856	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	26.69	276	373949	14.343	ng/ul#	74
25) Dibenzo(a,h)anthracene	26.67	278	94133	4.285	ng/ul	97
26) Benzo(a,h,i)perylene	27.53	276	307385	14.683	ng/ul	96

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

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