

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE102517\
 Data File : BE094548.D
 Acq On : 26 Oct 2017 00:24
 Operator : SJ/JU
 Sample : I5747-01 5X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-23(0-1)-100517

Manual Integrations
 APPROVED

Sohil
 10/27/2017 8:56:33 AM

Quant Time: Oct 26 13:14:21 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 25 18:44:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1543	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	8815	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	5015	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	11346	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	11125	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	10576	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	514	0.04	ng/ul	-0.01
14) Fluoranthene-d10	19.28	212	1297	0.04	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.74	128	2418	0.112	ng/ul#	93
5) 2-Methylnaphthalene	12.35	142	1231	0.082	ng/ul	95
7) Acenaphthylene	14.24	152	3788	0.167	ng/ul	99
8) Acenaphthene	14.58	153	3220	0.198	ng/ul	98
9) Fluorene	15.57	166	3063	0.167	ng/ul	94
12) Phenanthrene	17.29	178	68550	2.287	ng/ul#	89
13) Anthracene	17.38	178	14744	0.499	ng/ul#	90
15) Fluoranthene	19.31	202	154053	4.446	ng/ul	84
17) Pyrene	19.67	202	138075	4.015	ng/ul#	80
18) Benzo(a)anthracene	21.40	228	74580	2.345	ng/ul#	86
19) Chrysene	21.46	228	78114	2.376	ng/ul	99
21) Benzo(b)fluoranthene	23.18	252	109383m	3.677	ng/ul	
22) Benzo(k)fluoranthene	23.22	252	37676m	1.182	ng/ul	
23) Benzo(a)pyrene	23.84	252	72547	2.440	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	26.65	276	53813	1.749	ng/ul#	75
25) Dibenzo(a,h)anthracene	26.64	278	12673	0.489	ng/ul#	88
26) Benzo(g,h,i)perylene	27.48	276	44352	1.795	ng/ul	98

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

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