

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080818\
 Data File : BE097237.D
 Acq On : 8 Aug 2018 16:19
 Operator : SJ/JU
 Sample : J4268-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AA5

Manual Integrations
 APPROVED

Sohil
 8/9/2018 4:43:13 PM

Quant Time: Aug 09 01:03:19 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE080118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 08 23:25:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1255	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	6747	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	4020	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	9234	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	9655	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	8931m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	1580	0.15	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	4316	0.14	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.18	128	794	0.049	ng/ul#	85
7) Acenaphthylene	13.73	152	4532	0.261	ng/ul	98
8) Acenaphthene	14.08	153	1788	0.142	ng/ul	94
9) Fluorene	15.08	166	4203	0.291	ng/ul	93
12) Phenanthrene	16.81	178	26707	1.135	ng/ul#	89
13) Anthracene	16.90	178	6476	0.295	ng/ul#	92
15) Fluoranthene	18.84	202	65915	2.118	ng/ul	84
17) Pyrene	19.21	202	55153	2.132	ng/ul#	78
18) Benzo(a)anthracene	20.96	228	30425	1.103	ng/ul#	84
19) Chrysene	21.01	228	29266	1.029	ng/ul	98
21) Benzo(b)fluoranthene	22.53	252	40883m	1.680	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	11309m	0.431	ng/ul	
23) Benzo(a)pyrene	23.11	252	29031	1.192	ng/ul#	77
24) Indeno(1,2,3-cd)pyrene	25.50	276	19125	0.627	ng/ul#	71
25) Dibenzo(a,h)anthracene	25.52	278	4793	0.192	ng/ul	100
26) Benzo(a,h,i)perylene	26.20	276	17994	0.683	ng/ul#	93

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

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