

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022520\
 Data File : BM025100.D
 Acq On : 27 Feb 2020 04:53
 Operator : CG/JU
 Sample : L1506-20
 Misc :
 ALS Vial : 64 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBCZ8

Manual Integrations
 APPROVED

mohammad
 2/27/2020 5:01:43 PM

Quant Time: Feb 27 06:48:29 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM022420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 27 05:28:30 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	3048	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	11667	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.32	164	7671	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	17243m	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	18076	0.40	ng/ul	-0.01
20) Perylene-d12	23.44	264	21039	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.06	152	7197	0.33	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	19527	0.33	ng/ul	0.00
Target Compounds				Ovalue		
8) Acenaphthene	14.38	153	894	0.030	ng/ul	97
9) Fluorene	15.37	166	1026	0.028	ng/ul	99
12) Phenanthrene	17.10	178	1629m	0.027	ng/ul	
13) Anthracene	17.19	178	3214	0.055	ng/ul	97
15) Fluoranthene	19.12	202	7966	0.090	ng/ul	97
17) Pyrene	19.48	202	10913m	0.134	ng/ul	
18) Benzo(a)anthracene	21.23	228	4804	0.062	ng/ul	93
19) Chrysene	21.26	228	7987	0.103	ng/ul	98
21) Benzo(b)fluoranthene	22.79	252	19628m	0.238	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	8652m	0.106	ng/ul	
23) Benzo(a)pyrene	23.35	252	6551	0.087	ng/ul#	71
24) Indeno(1,2,3-cd)pyrene	25.63	276	7597	0.081	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.63	278	2048	0.026	ng/ul#	15
26) Benzo(g,h,i)perylene	26.29	276	5615	0.071	ng/ul	96

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

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