

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022520\
 Data File : BM025096.D
 Acq On : 27 Feb 2020 01:53
 Operator : CG/JU
 Sample : L1506-16
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBCZ4

Manual Integrations
 APPROVED

mohammad
 2/27/2020 5:01:33 PM

Quant Time: Feb 27 06:17:32 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	3066	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	12368	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.32	164	8256	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	18724m	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	20010	0.40	ng/ul	-0.01
20) Perylene-d12	23.44	264	22868	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.06	152	8501	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	23548	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.51	128	1386	0.037	ng/ul	98
7) Acenaphthylene	14.03	152	1342	0.028	ng/ul#	85
8) Acenaphthene	14.38	153	10307	0.322	ng/ul	99
9) Fluorene	15.37	166	13638	0.340	ng/ul	98
11) Pentachlorophenol	16.71	266	373	0.065	ng/ul	100
12) Phenanthrene	17.10	178	18326	0.278	ng/ul	96
13) Anthracene	17.19	178	3862	0.061	ng/ul	97
15) Fluoranthene	19.12	202	53551	0.560	ng/ul	96
17) Pyrene	19.48	202	45900m	0.509	ng/ul	
18) Benzo(a)anthracene	21.23	228	8214	0.095	ng/ul	99
19) Chrysene	21.28	228	8897m	0.103	ng/ul	
21) Benzo(b)fluoranthene	22.79	252	7143m	0.080	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	2667m	0.030	ng/ul	
23) Benzo(a)pyrene	23.35	252	2086	0.025	ng/ul#	64

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

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