

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022520\
 Data File : BM025090.D
 Acq On : 26 Feb 2020 22:17
 Operator : CG/JU
 Sample : L1506-10
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBCY8

Manual Integrations
 APPROVED

mohammad
 2/27/2020 5:01:14 PM

Quant Time: Feb 27 05:52:23 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	2540	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	10013	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.32	164	6853	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	16189m	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	17945	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	20717	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	5864	0.31	ng/ul	-0.01
14) Fluoranthene-d10	19.09	212	18009	0.32	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.10	178	1532	0.027	ng/ul#	85
15) Fluoranthene	19.12	202	7710	0.093	ng/ul	99
17) Pyrene	19.48	202	6723	0.083	ng/ul	97
18) Benzo(a)anthracene	21.23	228	1912m	0.025	ng/ul	
19) Chrysene	21.28	228	4835	0.063	ng/ul	96
21) Benzo(b)fluoranthene	22.79	252	6839m	0.084	ng/ul	
22) Benzo(k)fluoranthene	22.83	252	2310m	0.029	ng/ul	
23) Benzo(a)pyrene	23.35	252	2063	0.028	ng/ul#	64
24) Indeno(1,2,3-cd)pyrene	25.63	276	2253	0.024	ng/ul#	100
26) Benzo(g,h,i)perylene	26.30	276	2065	0.026	ng/ul#	74

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

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