

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022520\
 Data File : BM025094.D
 Acq On : 27 Feb 2020 00:41
 Operator : CG/JU
 Sample : L1506-14
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBCZ2

Manual Integrations
 APPROVED

mohammad
 2/27/2020 5:01:26 PM

Quant Time: Feb 27 06:08:44 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	3067	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	11958	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.32	164	7764	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	17270m	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	17912	0.40	ng/ul	-0.01
20) Perylene-d12	23.44	264	20613	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	7181	0.32	ng/ul	-0.01
14) Fluoranthene-d10	19.09	212	18450	0.31	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.04	152	1124	0.025	ng/ul#	86
13) Anthracene	17.19	178	1217	0.021	ng/ul#	86
15) Fluoranthene	19.12	202	3486	0.040	ng/ul	94
17) Pyrene	19.48	202	4050	0.050	ng/ul#	91
18) Benzo(a)anthracene	21.23	228	2592	0.034	ng/ul	94
19) Chrysene	21.28	228	3797m	0.049	ng/ul	
21) Benzo(b)fluoranthene	22.79	252	8018m	0.099	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	2934m	0.037	ng/ul	
23) Benzo(a)pyrene	23.35	252	2113	0.029	ng/ul#	64
24) Indeno(1,2,3-cd)pyrene	25.63	276	2351	0.025	ng/ul#	100
26) Benzo(g,h,i)perylene	26.30	276	1796	0.023	ng/ul#	76

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

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