

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025291.D
 Acq On : 05 Mar 2020 17:07
 Operator : CG/JU
 Sample : L1543-20
 Misc :
 ALS Vial : 87 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDA2

Manual Integrations
 APPROVED

mohammad
 3/6/2020 10:14:50 PM

Quant Time: Mar 06 02:52:22 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 04 08:02:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	3168	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	13116	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	9566	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	23423	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	21009	0.40	ng/ul	0.00
20) Perylene-d12	23.42	264	18602	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	1917	0.08	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	5960	0.08	ng/ul	0.00
Target Compounds				Ovalue		
13) Anthracene	17.17	178	3768	0.056	ng/ul	98
15) Fluoranthene	19.10	202	2832	0.030	ng/ul	99
17) Pyrene	19.46	202	3602m	0.045	ng/ul	
18) Benzo(a)anthracene	21.21	228	1692	0.022	ng/ul	94
19) Chrysene	21.27	228	4574m	0.054	ng/ul	
21) Benzo(b)fluoranthene	22.77	252	12494m	0.168	ng/ul	
22) Benzo(k)fluoranthene	22.81	252	3696m	0.047	ng/ul	
23) Benzo(a)pyrene	23.32	252	1826	0.029	ng/ul#	52
24) Indeno(1,2,3-cd)pyrene	25.60	276	2990	0.038	ng/ul#	100
26) Benzo(g,h,i)perylene	26.26	276	1490	0.023	ng/ul#	86

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

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