

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061120\
 Data File : BM026351.D
 Acq On : 11 Jun 2020 12:04
 Operator : JU/CG
 Sample : L2965-04
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SIDEWALK-COMP

Manual Integrations
 APPROVED

mohammad
 6/15/2020 8:43:14 AM

Quant Time: Jun 11 12:51:49 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM060520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 09 14:46:47 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.42	152	134868	20.00	ng	-0.01
21) Naphthalene-d8	10.17	136	528365	20.00	ng	-0.01
39) Acenaphthene-d10	14.07	164	300674	20.00	ng	-0.01
64) Phenanthrene-d10	16.82	188	630356	20.00	ng	-0.01
76) Chrysene-d12	21.03	240	696064	20.00	ng	-0.01
86) Perylene-d12	23.13	264	785797	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.05	112	392980	51.52	ng	-0.01
7) Phenol-d6	6.62	99	677811	61.36	ng	-0.01
23) Nitrobenzene-d5	8.56	82	522770	48.58	ng	-0.01
42) 2,4,6-Tribromophenol	15.57	330	215599	59.19	ng	-0.02
45) 2-Fluorobiphenyl	12.68	172	1028656	51.93	ng	-0.01
79) Terphenyl-d14	19.48	244	1657078	46.21	ng	-0.01
Target Compounds						
50) Dimethylphthalate	13.54	163	95227	4.330	ng	97
71) Phenanthrene	16.86	178	68205	2.078	ng	98
75) Fluoranthene	18.89	202	211872	5.630	ng	99
78) Pyrene	19.25	202	170906	3.716	ng	100
83) Chrysene	21.07	228	103585	2.602	ng	97
88) Benzo(b)fluoranthene	22.52	252	157411m	3.330	ng	
90) Benzo(a)pyrene	23.04	252	87075	2.074	ng	95
92) Benzo(g,h,i)perylene	25.79	276	93514	2.590	ng	98

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

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