

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091420\
 Data File : BP003279.D
 Acq On : 14 Sep 2020 21:13
 Operator : CG/JU
 Sample : L3981-11
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 B20-4-6

Manual Integrations
 APPROVED

mohammad
 9/15/2020 1:22:04 PM

Quant Time: Sep 22 08:44:30 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP082420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 01 14:37:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	201910	20.00	ng	-0.01
21) Naphthalene-d8	10.42	136	806066	20.00	ng	-0.01
39) Acenaphthene-d10	14.29	164	486612	20.00	ng	0.00
64) Phenanthrene-d10	17.04	188	985771	20.00	ng	0.00
76) Chrysene-d12	21.13	240	775722	20.00	ng	-0.01
86) Perylene-d12	23.37	264	841676	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	870449	76.22	ng	0.00
7) Phenol-d6	6.83	99	1092784	76.33	ng	0.00
23) Nitrobenzene-d5	8.79	82	664261	59.41	ng	-0.01
42) 2,4,6-Tribromophenol	15.79	330	546203	82.99	ng	0.00
45) 2-Fluorobiphenyl	12.90	172	1780221	53.58	ng	-0.01
79) Terphenyl-d14	19.62	244	2129278	60.40	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.75	163	239946	6.466	ng	99
75) Fluoranthene	19.06	202	382654	6.198	ng	100
78) Pyrene	19.41	202	342521	6.803	ng	99
81) Benzo(a)anthracene	21.12	228	212546	4.374	ng	96
83) Chrysene	21.17	228	215494	4.632	ng	97
88) Benzo(b)fluoranthene	22.70	252	302014m	5.773	ng	
90) Benzo(a)pyrene	23.27	252	199174	4.104	ng	97
92) Benzo(g,h,i)perylene	26.33	276	151061	2.919	ng	98

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

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