

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091420\
 Data File : BP003277.D
 Acq On : 14 Sep 2020 20:05
 Operator : CG/JU
 Sample : L3993-05
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PAV-091120-DP

Manual Integrations
 APPROVED

mohammad
 9/15/2020 1:21:10 PM

Quant Time: Sep 23 07:25:28 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	193873	20.00	ng	-0.01
21) Naphthalene-d8	10.42	136	779433	20.00	ng	-0.02
39) Acenaphthene-d10	14.29	164	469616	20.00	ng	0.00
64) Phenanthrene-d10	17.04	188	972462	20.00	ng	0.00
76) Chrysene-d12	21.13	240	810299	20.00	ng	-0.01
86) Perylene-d12	23.36	264	865321	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	875132	79.81	ng	0.00
7) Phenol-d6	6.83	99	1053871	76.67	ng	0.00
23) Nitrobenzene-d5	8.79	82	656769	60.75	ng	-0.01
42) 2,4,6-Tribromophenol	15.78	330	558628	87.95	ng	-0.01
45) 2-Fluorobiphenyl	12.90	172	1779222	55.48	ng	-0.01
79) Terphenyl-d14	19.62	244	2271538	61.69	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.75	163	222310	6.208	ng	99
71) Phenanthrene	17.08	178	179003	3.397	ng	99
75) Fluoranthene	19.06	202	405261	6.654	ng	99
78) Pyrene	19.41	202	380909	7.243	ng	99
81) Benzo(a)anthracene	21.12	228	215055	4.237	ng	99
83) Chrysene	21.17	228	192981	3.971	ng	97
88) Benzo(b)fluoranthene	22.70	252	265449m	4.936	ng	
90) Benzo(a)pyrene	23.27	252	201783	4.044	ng	95
92) Benzo(g,h,i)perylene	26.33	276	128349	2.413	ng	97

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

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