

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP072120\
 Data File : BP002817.D
 Acq On : 21 Jul 2020 16:54
 Operator : CG/JU
 Sample : L3360-03
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SP-2

Manual Integrations
 APPROVED

Sohil
 7/24/2020 12:10:24 PM

Quant Time: Jul 21 17:54:59 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 13 10:54:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	122273	20.00	ng	0.00
21) Naphthalene-d8	10.32	136	435236	20.00	ng	-0.01
39) Acenaphthene-d10	14.20	164	258058	20.00	ng	-0.01
64) Phenanthrene-d10	16.96	188	583284	20.00	ng	-0.01
76) Chrysene-d12	21.07	240	637812	20.00	ng	0.00
86) Perylene-d12	23.26	264	685850	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.18	112	133310	18.39	ng	0.00
7) Phenol-d6	6.75	99	510738	49.38	ng	0.00
23) Nitrobenzene-d5	8.69	82	438453	53.88	ng	-0.01
42) 2,4,6-Tribromophenol	15.70	330	49185	18.26	ng	0.00
45) 2-Fluorobiphenyl	12.81	172	886820	53.16	ng	-0.01
79) Terphenyl-d14	19.55	244	1366855	49.13	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.66	163	138653	7.045	ng	99
71) Phenanthrene	17.00	178	73887	2.340	ng	99
75) Fluoranthene	18.99	202	142603	3.897	ng	97
78) Pyrene	19.34	202	144495	3.281	ng	99
81) Benzo(a)anthracene	21.06	228	86670	2.099	ng	94
83) Chrysene	21.10	228	90596m	2.305	ng	
88) Benzo(b)fluoranthene	22.61	252	112854m	2.650	ng	
90) Benzo(a)pyrene	23.17	252	82406	2.042	ng	# 83

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

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