

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070220\
 Data File : BP002638.D
 Acq On : 02 Jul 2020 19:21
 Operator : CG/JU
 Sample : L3188-01 5X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TR-04-070120

Manual Integrations
 APPROVED

mohammad
 7/7/2020 8:25:02 AM

Quant Time: Jul 02 19:54:50 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 29 16:51:57 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	169124	20.00	ng	0.00
21) Naphthalene-d8	10.34	136	707421	20.00	ng	0.00
39) Acenaphthene-d10	14.22	164	463105	20.00	ng	0.00
64) Phenanthrene-d10	16.97	188	1067760	20.00	ng	0.00
76) Chrysene-d12	21.09	240	1135453	20.00	ng	0.00
86) Perylene-d12	23.28	264	1052989	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.19	112	121796	12.19	ng	0.00
7) Phenol-d6	6.76	99	181494	13.02	ng	0.00
23) Nitrobenzene-d5	8.72	82	122930	8.95	ng	0.00
42) 2,4,6-Tribromophenol	15.72	330	66746	11.85	ng	0.00
45) 2-Fluorobiphenyl	12.83	172	317827	10.10	ng	0.00
79) Terphenyl-d14	19.57	244	547507	10.41	ng	0.00
Target Compounds						
71) Phenanthrene	17.02	178	180948	3.150	ng	99
75) Fluoranthene	19.01	202	311893	4.485	ng	98
78) Pyrene	19.36	202	345021	4.464	ng	98
81) Benzo(a)anthracene	21.07	228	212309	2.860	ng	99
83) Chrysene	21.12	228	201786	2.835	ng	97
88) Benzo(b)fluoranthene	22.63	252	250260m	3.725	ng	
90) Benzo(a)pyrene	23.19	252	192404	3.120	ng	# 96

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

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