

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP073020\
 Data File : BP002879.D
 Acq On : 30 Jul 2020 16:42
 Operator : CG/JU
 Sample : L3473-09
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-2-D

Quant Time: Jul 30 17:41:33 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 30 15:30:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	512506	20.00	ng	0.00
21) Naphthalene-d8	10.50	136	1999225	20.00	ng	0.00
39) Acenaphthene-d10	14.36	164	1237470	20.00	ng	0.00
64) Phenanthrene-d10	17.11	188	2605368	20.00	ng	0.00
76) Chrysene-d12	21.20	240	2686986	20.00	ng	0.00
86) Perylene-d12	23.47	264	2850205	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	2615144	78.10	ng	0.00
7) Phenol-d6	6.89	99	3351595	72.09	ng	0.00
23) Nitrobenzene-d5	8.86	82	2074804	56.21	ng	0.00
42) 2,4,6-Tribromophenol	15.85	330	1253092	81.47	ng	0.00
45) 2-Fluorobiphenyl	12.98	172	4670563	51.57	ng	0.00
79) Terphenyl-d14	19.69	244	6364839	47.42	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	13.82	163	730941	7.552	ng	100
71) Phenanthrene	17.15	178	381049	2.494	ng	98
75) Fluoranthene	19.13	202	778144	4.421	ng	98
78) Pyrene	19.47	202	695828	3.704	ng	98
81) Benzo(a)anthracene	21.19	228	445286	2.433	ng	99
83) Chrysene	21.24	228	357837	2.073	ng	97
88) Benzo(b)fluoranthene	22.79	252	514817	2.560	ng	# 98
90) Benzo(a)pyrene	23.37	252	375070	2.064	ng	98

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

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