

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070921\
 Data File : BP006166.D
 Acq On : 09 Jul 2021 18:06
 Operator : CG/JU
 Sample : M2969-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB01

Quant Time: Jul 10 01:12:02 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP070721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 07 16:08:13 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	211107	20.000	ng	0.00
21) Naphthalene-d8	10.675	136	794594	20.000	ng	0.00
39) Acenaphthene-d10	14.510	164	467681	20.000	ng	0.00
64) Phenanthrene-d10	17.263	188	908203	20.000	ng	# 0.00
76) Chrysene-d12	21.345	240	867761	20.000	ng	# 0.00
86) Perylene-d12	23.739	264	978242	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	1112455	90.671	ng	-0.01
7) Phenol-d6	7.058	99	1311704	80.826	ng	0.00
23) Nitrobenzene-d5	9.040	82	742604	59.032	ng	0.00
42) 2,4,6-Tribromophenol	16.004	330	448909	80.622	ng	0.00
45) 2-Fluorobiphenyl	13.134	172	1575987	49.817	ng	0.00
79) Terphenyl-d14	19.816	244	2229822	50.663	ng	0.00
Target Compounds						Qvalue
75) Fluoranthene	19.275	202	144519	2.571	ng	96
78) Pyrene	19.628	202	143441	2.446	ng	98

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

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