

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070921\
 Data File : BP006176.D
 Acq On : 09 Jul 2021 23:48
 Operator : CG/JU
 Sample : M2969-13 5X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB07

Quant Time: Jul 10 01:14:43 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	220547	20.000	ng	0.00
21) Naphthalene-d8	10.675	136	858463	20.000	ng	0.00
39) Acenaphthene-d10	14.510	164	471554	20.000	ng	0.00
64) Phenanthrene-d10	17.263	188	856646	20.000	ng	# 0.00
76) Chrysene-d12	21.345	240	821734	20.000	ng	# 0.00
86) Perylene-d12	23.739	264	962189	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.458	112	97899	7.638	ng	0.00
7) Phenol-d6	7.063	99	280007	16.515	ng	0.00
23) Nitrobenzene-d5	9.046	82	154667	11.380	ng	0.00
42) 2,4,6-Tribromophenol	16.016	330	12893	2.297	ng	0.00
45) 2-Fluorobiphenyl	13.134	172	371263	11.639	ng	0.00
79) Terphenyl-d14	19.810	244	462955	11.108	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.304	178	97951	2.025	ng	99
75) Fluoranthene	19.269	202	128230	2.418	ng	94
78) Pyrene	19.622	202	118842	2.140	ng	98

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

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