

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060821\
 Data File : BP005805.D
 Acq On : 09 Jun 2021 10:25
 Operator : CG/JU
 Sample : M2589-30
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FOUNDATION-EXCAVATION-SAMP

Quant Time: Jun 09 12:24:59 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 09 11:54:55 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.957	152	74854	20.000	ng	0.00
21) Naphthalene-d8	10.763	136	271070	20.000	ng	0.00
39) Acenaphthene-d10	14.581	164	162308	20.000	ng	0.00
64) Phenanthrene-d10	17.327	188	340002	20.000	ng	0.00
76) Chrysene-d12	21.392	240	393370	20.000	ng	0.00
86) Perylene-d12	23.815	264	457995	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	571089	122.423	ng	0.00
7) Phenol-d6	7.122	99	583448	96.496	ng	0.00
23) Nitrobenzene-d5	9.122	82	476610	86.202	ng	0.00
42) 2,4,6-Tribromophenol	16.069	330	358883	128.848	ng	0.00
45) 2-Fluorobiphenyl	13.210	172	1033732	83.633	ng	0.00
79) Terphenyl-d14	19.869	244	1742621	82.951	ng	0.00

Target Compounds	Qvalue

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

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