

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060222\
 Data File : BP010666.D
 Acq On : 02 Jun 2022 12:16
 Operator : CG/JU
 Sample : N2953-10
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 P044-SS001-1824-01

Quant Time: Jun 03 15:43:09 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat May 28 02:18:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	108945	20.000	ng	-0.01
21) Naphthalene-d8	10.657	136	437439	20.000	ng	-0.01
39) Acenaphthene-d10	14.492	164	268386	20.000	ng	-0.01
64) Phenanthrene-d10	17.251	188	550508	20.000	ng	0.00
76) Chrysene-d12	21.339	240	583026	20.000	ng	# 0.00
86) Perylene-d12	23.745	264	598254	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	841538	140.917	ng	0.00
7) Phenol-d6	7.034	99	1128118	133.065	ng	0.00
23) Nitrobenzene-d5	9.016	82	745198	80.541	ng	-0.01
42) 2,4,6-Tribromophenol	15.992	330	483242	159.065	ng	0.00
45) 2-Fluorobiphenyl	13.116	172	1411224	74.626	ng	-0.01
79) Terphenyl-d14	19.809	244	2237630	72.089	ng	0.00

Target Compounds	Qvalue

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

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