

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060222\
 Data File : BP010667.D
 Acq On : 02 Jun 2022 12:50
 Operator : CG/JU
 Sample : N2953-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 P043-SS001-2430-01

Quant Time: Jun 03 15:43:37 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	128104	20.000	ng	-0.01
21) Naphthalene-d8	10.657	136	513765	20.000	ng	-0.01
39) Acenaphthene-d10	14.492	164	314820	20.000	ng	-0.01
64) Phenanthrene-d10	17.251	188	622439	20.000	ng	0.00
76) Chrysene-d12	21.333	240	591483	20.000	ng	0.00
86) Perylene-d12	23.745	264	624007	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	774177	110.249	ng	0.00
7) Phenol-d6	7.034	99	1026498	102.970	ng	0.00
23) Nitrobenzene-d5	9.022	82	698836	64.309	ng	0.00
42) 2,4,6-Tribromophenol	15.986	330	409001	114.771	ng	0.00
45) 2-Fluorobiphenyl	13.116	172	1247889	56.256	ng	-0.01
79) Terphenyl-d14	19.809	244	1781968	56.588	ng	0.00

Target Compounds	Qvalue

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

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