

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060322\
 Data File : BP010699.D
 Acq On : 03 Jun 2022 18:47
 Operator : CG/JU
 Sample : N3143-06
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 20220427-SPLP-AMENDED-SB-8-(DISCRETE-COMP-G)

Quant Time: Jun 04 02:01:38 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jun 04 01:56:43 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.852	152	109072	20.000	ng	0.00
21) Naphthalene-d8	10.651	136	428040	20.000	ng	0.00
39) Acenaphthene-d10	14.487	164	276564	20.000	ng	-0.01
64) Phenanthrene-d10	17.245	188	581597	20.000	ng	0.00
76) Chrysene-d12	21.333	240	589004	20.000	ng	#-0.01
86) Perylene-d12	23.739	264	605866	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	443164	74.122	ng	0.00
7) Phenol-d6	7.034	99	362724	42.734	ng	0.00
23) Nitrobenzene-d5	9.010	82	807264	89.165	ng	-0.01
42) 2,4,6-Tribromophenol	15.986	330	606820	193.836	ng	0.00
45) 2-Fluorobiphenyl	13.110	172	1785588	91.631	ng	0.00
79) Terphenyl-d14	19.804	244	3193797	101.849	ng	0.00

Target Compounds	Qvalue

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

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