

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040624\
 Data File : BG060811.D
 Acq On : 6 Apr 2024 14:52
 Operator : MA/JU
 Sample : P1995-06
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MOO-23-00113

Quant Time: Apr 08 02:06:46 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG032924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 30 03:12:26 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	68049	20.000	ng	0.00
21) Naphthalene-d8	10.749	136	259243	20.000	ng	0.00
39) Acenaphthene-d10	14.580	164	193182	20.000	ng	0.00
64) Phenanthrene-d10	17.330	188	432679	20.000	ng	0.00
76) Chrysene-d12	21.589	240	389256	20.000	ng	-0.01
86) Perylene-d12	24.727	264	471653	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.543	112	512015	125.345	ng	0.00
7) Phenol-d6	7.118	99	660530	108.936	ng	0.00
23) Nitrobenzene-d5	9.110	82	511776	112.646	ng	0.00
42) 2,4,6-Tribromophenol	16.072	330	388395	177.112	ng	0.00
45) 2-Fluorobiphenyl	13.211	172	1219040	92.612	ng	0.00
79) Terphenyl-d14	19.956	244	1932654	87.425	ng	0.00
Target Compounds						
32) Benzoic acid	10.003	122	7345	9.088	ng	Qvalue # 79

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

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