

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040624\
 Data File : BG060810.D
 Acq On : 6 Apr 2024 14:11
 Operator : MA/JU
 Sample : P1995-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MOO-23-00119

Quant Time: Apr 08 02:06:12 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	7.954	152	63693	20.000	ng	0.00
21) Naphthalene-d8	10.750	136	251066	20.000	ng	0.00
39) Acenaphthene-d10	14.581	164	194429	20.000	ng	0.00
64) Phenanthrene-d10	17.325	188	400759	20.000	ng	-0.01
76) Chrysene-d12	21.590	240	389206	20.000	ng	-0.01
86) Perylene-d12	24.728	264	449790	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.545	112	484003	126.591	ng	0.00
7) Phenol-d6	7.119	99	617187	108.749	ng	0.00
23) Nitrobenzene-d5	9.105	82	472391	107.364	ng	0.00
42) 2,4,6-Tribromophenol	16.068	330	364202	165.015	ng	0.00
45) 2-Fluorobiphenyl	13.206	172	1126997	85.070	ng	-0.01
79) Terphenyl-d14	19.957	244	1966477	88.967	ng	0.00
Target Compounds						
22) Acetophenone	8.764	105	16266	2.350	ng	Qvalue # 90
32) Benzoic acid	9.992	122	28383	15.823	ng	91

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

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