Quantitation Report (QT/LSC Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022322\
Data File : BM034053.D
Acq On : 23 Feb 2022 15:24
Operator : CG/JU
Sample : N1632-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
20220090-BERKELEY-2-ELUTRIATE

Quant Time: Feb 23 16:57:45 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM022222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 22 17:55:56 2022
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.019	152	248772	20.000	ng	0.00
21) Naphthalene-d8	10.837	136	936112	20.000	ng	# 0.00
39) Acenaphthene-d10	14.654	164	591780	20.000	ng	0.00
64) Phenanthrene-d10	17.395	188	1216251	20.000	ng	# 0.00
76) Chrysene-d12	21.565	240	1270426	20.000	ng	0.00
86) Perylene-d12	24.030	264	1360175	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.566	112	1009067	72.307	ng	0.00
7) Phenol-d6	7.166	99	878817	47.956	ng	0.00
23) Nitrobenzene-d5	9.178	82	1612665	92.968	ng	0.00
42) 2,4,6-Tribromophenol	16.136	330	1220048	169.869	ng	0.00
45) 2-Fluorobiphenyl	13.283	172	4182711	94.167	ng	0.00
79) Terphenyl-d14	20.012	244	5176818	70.957	ng	0.00
Target Compounds						Qvalue
10) Phenol	7.190	94	56152	3.115	ng	92

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

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