

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN010219\
 Data File : VN053246.D
 Acq On : 2 Jan 2019 16:01
 Operator : MD\SY
 Sample : K1023-01DL 10X
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 12 Sample Multiplier: 28

Instrument :
 MSVOA_N
 ClientSampleId :
 PURGE-WATERDL

Quant Time: Jan 03 09:55:09 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N121818W.M
 Quant Title : SW846 8260
 QLast Update : Tue Dec 18 13:03:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	1125329	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	1722854	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	1612665	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	606134	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	545719	47.79	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.58%	
35) Dibromofluoromethane	7.59	113	492105	62.53	ug/l	0.00
Spiked Amount 50.000			Recovery	=	125.06%	
50) Toluene-d8	10.09	98	2144935	50.77	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.54%	
62) 4-Bromofluorobenzene	12.40	95	734130	49.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.00%	
Target Compounds						
16) Acetone	3.82	43	12356	5.518	ug/l	98
27) cis-1,2-Dichloroethene	6.83	96	42057	3.224	ug/l	89
44) Trichloroethene	8.83	130	14919	1.049	ug/l	97
64) Tetrachloroethene	10.63	164	787656	45.089	ug/l	98

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

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