

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN042919\
 Data File : VN055293.D
 Acq On : 29 Apr 2019 15:03
 Operator : JC/SP
 Sample : K2598-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 OIL-WATER-SEPERATOR-2(OWS-2)

Quant Time: Apr 30 05:26:42 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N042919W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 30 04:20:38 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	377215	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	648277	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	592737	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	186808	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	8.03	65	261194	52.93	ug/l	0.00
Spiked Amount			Recovery	=	105.86%	
35) Dibromofluoromethane	7.59	113	212106	49.85	ug/l	0.00
Spiked Amount			Recovery	=	99.70%	
50) Toluene-d8	10.09	98	805267	52.41	ug/l	0.00
Spiked Amount			Recovery	=	104.82%	
62) 4-Bromofluorobenzene	12.40	95	253724	52.84	ug/l	0.00
Spiked Amount			Recovery	=	105.68%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
16) Acetone	3.82	43	81487	52.995	ug/l	100
17) Carbon Disulfide	4.05	76	12076	1.242	ug/l	97
18) Methyl Acetate	4.33	43	7267	0.924	ug/l #	67
20) Methylene Chloride	4.55	84	3180968	677.762	ug/l	98
25) 2-Butanone	6.84	43	32519	16.537	ug/l	94
43) Isopropyl Acetate	8.16	43	14825	1.938	ug/l #	90
52) Toluene	10.16	92	126620	12.056	ug/l	97
67) Ethyl Benzene	11.51	91	9009	0.430	ug/l	92
68) m/p-Xylenes	11.62	106	8681	1.162	ug/l	89
69) o-Xylene	11.95	106	6730	0.909	ug/l	97
78) n-propylbenzene	12.59	91	5738	0.299	ug/l	95
80) 1,3,5-Trimethylbenzene	12.73	105	11686	0.798	ug/l	99
84) 1,2,4-Trimethylbenzene	13.04	105	42399	3.237	ug/l	98
86) p-Isopropyltoluene	13.29	119	59720	4.565	ug/l	98

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

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