

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082119\
 Data File : BM022237.D
 Acq On : 21 Aug 2019 20:02
 Operator : HP/JU
 Sample : K4402-01RE
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 FRAC-TANK-MUDRE

Quant Time: Aug 22 03:45:21 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 20 02:18:05 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	44368	20.00	ng	-0.01
21) Naphthalene-d8	10.34	136	174039	20.00	ng	0.00
39) Acenaphthene-d10	14.22	164	101269	20.00	ng	0.00
64) Phenanthrene-d10	16.99	188	182663	20.00	ng	0.00
76) Chrysene-d12	21.19	240	161427	20.00	ng	0.00
87) Perylene-d12	23.39	264	193272	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.19	112	8677	3.48	ng	0.00
7) Phenol-d6	6.76	99	76309	23.01	ng	0.00
23) Nitrobenzene-d5	8.73	82	211648	59.15	ng	-0.01
42) 2,4,6-Tribromophenol	15.75	330	1537	0.94	ng	0.01
45) 2-Fluorobiphenyl	12.83	172	324688	40.05	ng	-0.01
79) Terphenyl-d14	19.63	244	260489	23.51	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.70	163	44654	5.245	ng	Qvalue # 99

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

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