

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110421\
 Data File : BM032880.D
 Acq On : 04 Nov 2021 18:12
 Operator : CG/JU
 Sample : M4499-02
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WEL-BRIDGE

Quant Time: Nov 07 08:50:21 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM110221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 02 16:07:18 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.484	152	105158	20.000	ng	0.00
21) Naphthalene-d8	10.266	136	405806	20.000	ng	# 0.00
39) Acenaphthene-d10	14.148	164	232826	20.000	ng	0.00
64) Phenanthrene-d10	16.883	188	445682	20.000	ng	-0.01
76) Chrysene-d12	21.065	240	421055	20.000	ng	-0.01
86) Perylene-d12	23.177	264	444431	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.084	112	893833	148.187	ng	0.00
7) Phenol-d6	6.695	99	1042020	116.005	ng	0.00
23) Nitrobenzene-d5	8.642	82	725799	97.892	ng	0.00
42) 2,4,6-Tribromophenol	15.642	330	353074	135.407	ng	0.00
45) 2-Fluorobiphenyl	12.760	172	1534209	102.202	ng	0.00
79) Terphenyl-d14	19.518	244	1911461	99.113	ng	0.00

Target Compounds	Qvalue

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

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