

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121721\
 Data File : BM033506.D
 Acq On : 17 Dec 2021 15:12
 Operator : CG/JU
 Sample : M5064-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SASP2-121-1-(5-10)

Quant Time: Dec 18 04:51:17 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM121621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 16 17:10:15 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.937	152	52080	20.000	ng	0.00
21) Naphthalene-d8	10.748	136	211472	20.000	ng	# 0.00
39) Acenaphthene-d10	14.583	164	140650	20.000	ng	0.00
64) Phenanthrene-d10	17.330	188	272199	20.000	ng	0.00
76) Chrysene-d12	21.512	240	160538	20.000	ng	# 0.00
86) Perylene-d12	23.924	264	138744	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.502	112	334012	110.667	ng	0.00
7) Phenol-d6	7.113	99	440307	108.707	ng	0.00
23) Nitrobenzene-d5	9.113	82	341029	69.243	ng	0.00
42) 2,4,6-Tribromophenol	16.077	330	151455	95.727	ng	0.00
45) 2-Fluorobiphenyl	13.207	172	696434	68.495	ng	0.00
79) Terphenyl-d14	19.954	244	838426	82.221	ng	0.00

Target Compounds	Qvalue

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

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