

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090921\
 Data File : BM032110.D
 Acq On : 09 Sep 2021 13:51
 Operator : CG/JU
 Sample : M3644-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 191-0902-DUP

Quant Time: Sep 09 17:01:11 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM090821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 08 19:23:17 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.357	152	66468	20.000	ng	0.00
21) Naphthalene-d8	10.116	136	261286	20.000	ng	0.00
39) Acenaphthene-d10	14.015	164	173354	20.000	ng	0.00
64) Phenanthrene-d10	16.786	188	326268	20.000	ng	0.00
76) Chrysene-d12	21.003	240	248388	20.000	ng	0.00
86) Perylene-d12	23.097	264	245236	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.004	112	101349	27.208	ng	0.00
7) Phenol-d6	6.563	99	88359	15.592	ng	0.00
23) Nitrobenzene-d5	8.510	82	550174	93.841	ng	0.00
42) 2,4,6-Tribromophenol	15.527	330	65050	47.217	ng	0.00
45) 2-Fluorobiphenyl	12.621	172	1184427	89.584	ng	0.00
79) Terphenyl-d14	19.444	244	1415411	92.975	ng	0.00

Target Compounds	Qvalue

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

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