

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101321\
 Data File : BM032495.D
 Acq On : 13 Oct 2021 14:37
 Operator : CG/JU
 Sample : M4150-02
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 FDR-20211007-WC

Quant Time: Oct 13 16:40:59 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM100221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 13 12:02:28 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.601	152	230897	20.00	ng	0.00
21) Naphthalene-d8	10.389	136	885065	20.00	ng	# 0.00
39) Acenaphthene-d10	14.248	164	496317	20.00	ng	0.00
64) Phenanthrene-d10	16.977	188	916215	20.00	ng	# 0.00
76) Chrysene-d12	21.142	240	759046	20.00	ng	# 0.00
86) Perylene-d12	23.283	264	749495	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.184	112	1797474	131.17	ng	0.00
7) Phenol-d6	6.795	99	2090443	111.55	ng	0.00
23) Nitrobenzene-d5	8.754	82	1507013	94.70	ng	0.00
42) 2,4,6-Tribromophenol	15.736	330	691843	124.52	ng	0.00
45) 2-Fluorobiphenyl	12.866	172	3066940	91.88	ng	0.00
79) Terphenyl-d14	19.595	244	3463343	96.09	ng	0.00

Target Compounds	Qvalue

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

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