

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101321\
 Data File : BM032500.D
 Acq On : 13 Oct 2021 17:39
 Operator : CG/JU
 Sample : M4047-02
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 149-ENV-1-GRAB

Quant Time: Oct 14 02:31:36 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	229210	20.00	ng	0.00
21) Naphthalene-d8	10.389	136	889689	20.00	ng	# 0.00
39) Acenaphthene-d10	14.248	164	506833	20.00	ng	0.00
64) Phenanthrene-d10	16.977	188	945961	20.00	ng	# 0.00
76) Chrysene-d12	21.142	240	780901	20.00	ng	# 0.00
86) Perylene-d12	23.288	264	791047	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.184	112	1651181	121.38	ng	0.00
7) Phenol-d6	6.795	99	1813556	97.49	ng	0.00
23) Nitrobenzene-d5	8.754	82	1482115	92.65	ng	0.00
42) 2,4,6-Tribromophenol	15.736	330	715477	126.10	ng	0.00
45) 2-Fluorobiphenyl	12.866	172	2999061	87.98	ng	0.00
79) Terphenyl-d14	19.595	244	3580142	96.55	ng	0.00

Target Compounds	Qvalue

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

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