

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071824\
 Data File : BM046705.D
 Acq On : 18 Jul 2024 16:41
 Operator : MA/JU
 Sample : P3246-22
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-CTW2-BL-05

Quant Time: Jul 18 17:15:55 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM071724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 18 06:01:46 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.487	152	94245	20.000	ng	0.00
21) Naphthalene-d8	10.246	136	315611	20.000	ng	# 0.00
39) Acenaphthene-d10	14.133	164	211245	20.000	ng	0.00
64) Phenanthrene-d10	16.886	188	414796	20.000	ng	0.00
76) Chrysene-d12	21.115	240	301984	20.000	ng	0.00
86) Perylene-d12	23.880	264	344941	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.134	112	667844	143.933	ng	0.00
7) Phenol-d6	6.693	99	736981	124.735	ng	0.00
23) Nitrobenzene-d5	8.634	82	642735	105.734	ng	0.00
42) 2,4,6-Tribromophenol	15.633	330	472910	144.330	ng	0.00
45) 2-Fluorobiphenyl	12.751	172	1640126	109.670	ng	0.00
79) Terphenyl-d14	19.545	244	2151278	128.827	ng	0.00

Target Compounds	Qvalue

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

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