

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072122\
 Data File : BM035895.D
 Acq On : 21 Jul 2022 18:55
 Operator : CG/JU
 Sample : N3786-17
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 0718-WC-TP-WC17A

Quant Time: Jul 22 01:08:51 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM071822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 18 17:08:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.892	152	344002	20.000	ng	0.00
21) Naphthalene-d8	10.704	136	1455077	20.000	ng	0.00
39) Acenaphthene-d10	14.527	164	851650	20.000	ng	0.00
64) Phenanthrene-d10	17.268	188	1640254	20.000	ng	0.00
76) Chrysene-d12	21.444	240	1296519	20.000	ng	0.00
86) Perylene-d12	23.821	264	1241201	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.451	112	1056697	47.622	ng	0.00
7) Phenol-d6	7.057	99	1347414	48.289	ng	0.00
23) Nitrobenzene-d5	9.063	82	813579	29.904	ng	0.00
42) 2,4,6-Tribromophenol	16.010	330	437033	45.877	ng	-0.01
45) 2-Fluorobiphenyl	13.163	172	1850485	30.713	ng	0.00
79) Terphenyl-d14	19.892	244	2298713	34.496	ng	0.00

Target Compounds	Qvalue

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

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