

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM103122\
 Data File : BM037287.D
 Acq On : 31 Oct 2022 13:28
 Operator : CG/JU
 Sample : N5272-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 20221102-COMP-01

Quant Time: Nov 01 05:01:18 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 31 01:34:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	277770	20.000	ng	-0.01
21) Naphthalene-d8	10.669	136	1094691	20.000	ng	0.00
39) Acenaphthene-d10	14.504	164	601047	20.000	ng	0.00
64) Phenanthrene-d10	17.245	188	1183016	20.000	ng	0.00
76) Chrysene-d12	21.415	240	1284063	20.000	ng	0.00
86) Perylene-d12	23.780	264	1480657	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	1823215	113.398	ng	0.00
7) Phenol-d6	7.040	99	2252532	103.226	ng	0.00
23) Nitrobenzene-d5	9.034	82	1344722	61.976	ng	-0.01
42) 2,4,6-Tribromophenol	15.992	330	778962	109.974	ng	0.00
45) 2-Fluorobiphenyl	13.127	172	2634117	58.231	ng	-0.01
79) Terphenyl-d14	19.862	244	3593180	50.181	ng	0.00

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

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