

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM052522\
 Data File : BM035258.D
 Acq On : 26 May 2022 01:10
 Operator : CG/JU
 Sample : N2922-21
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P024-SS003-0006-01

Quant Time: May 26 05:55:00 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM042722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 26 05:05:10 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	83377	20.000	ng	0.00
21) Naphthalene-d8	10.687	136	301048	20.000	ng	# 0.00
39) Acenaphthene-d10	14.522	164	184002	20.000	ng	0.00
64) Phenanthrene-d10	17.268	188	398161	20.000	ng	# 0.00
76) Chrysene-d12	21.451	240	526904	20.000	ng	# 0.00
86) Perylene-d12	23.821	264	626087	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	580291	122.787	ng	0.00
7) Phenol-d6	7.057	99	708737	116.168	ng	0.00
23) Nitrobenzene-d5	9.045	82	434283	77.901	ng	0.00
42) 2,4,6-Tribromophenol	16.010	330	367621	183.666	ng	0.00
45) 2-Fluorobiphenyl	13.151	172	1149208	90.666	ng	0.00
79) Terphenyl-d14	19.892	244	1942405	71.377	ng	0.00
Target Compounds						
84) Bis(2-ethylhexyl)phtha...	21.368	149	210687	8.958	ng	# 96

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

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