

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080825\
 Data File : BP025332.D
 Acq On : 08 Aug 2025 14:30
 Operator : CG/JU
 Sample : Q2759-02DL 10X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SOLID-DRUMDL

Quant Time: Aug 08 14:49:51 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 06 06:41:44 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.801	152	317812	20.000	ng	0.00
21) Naphthalene-d8	10.578	136	1234127	20.000	ng	0.00
39) Acenaphthene-d10	14.431	164	717730	20.000	ng	0.01
64) Phenanthrene-d10	17.224	188	1200789	20.000	ng	0.00
76) Chrysene-d12	21.677	240	1071297	20.000	ng	0.02
86) Perylene-d12	25.065	264	1324881	20.000	ng	0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.407	112	206586	10.455	ng	0.00
7) Phenol-d6	6.960	99	235667	9.062	ng	0.00
23) Nitrobenzene-d5	8.931	82	173793	7.802	ng	-0.01
42) 2,4,6-Tribromophenol	15.930	330	83741	11.695	ng	0.00
45) 2-Fluorobiphenyl	13.042	172	396530	7.631	ng	0.01
79) Terphenyl-d14	19.948	244	390132	6.878	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	6.990	94	75364	2.709	ng	96
20) 3+4-Methylphenols	8.543	107	671730	26.201	ng	90

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

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