

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040925\
 Data File : BP024237.D
 Acq On : 09 Apr 2025 16:20
 Operator : RC/JU
 Sample : Q1712-08
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TT-7

Quant Time: Apr 09 16:47:54 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031225.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 13 05:55:58 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.740	152	256391	20.000	ng	-0.03
21) Naphthalene-d8	10.516	136	980816	20.000	ng	-0.04
39) Acenaphthene-d10	14.375	164	587798	20.000	ng	-0.03
64) Phenanthrene-d10	17.186	188	1203665	20.000	ng	-0.02
76) Chrysene-d12	21.621	240	1296452	20.000	ng	-0.04
86) Perylene-d12	24.998	264	1347892	20.000	ng	-0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.358	112	2113496	131.586	ng	-0.02
7) Phenol-d6	6.922	99	2478651	115.400	ng	-0.03
23) Nitrobenzene-d5	8.887	82	1684087	96.876	ng	-0.04
42) 2,4,6-Tribromophenol	15.892	330	1168202	157.778	ng	-0.02
45) 2-Fluorobiphenyl	12.987	172	3616955	90.733	ng	-0.03
79) Terphenyl-d14	19.898	244	6180318	98.404	ng	-0.02

Target Compounds	Qvalue

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

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