

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061125\
 Data File : BP024914.D
 Acq On : 11 Jun 2025 16:24
 Operator : RC/JU
 Sample : Q2244-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP03-MHC

Quant Time: Jun 11 16:59:16 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060625.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 06 16:20:27 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.607	152	248278	20.000	ng	0.00
21) Naphthalene-d8	10.384	136	971478	20.000	ng	0.00
39) Acenaphthene-d10	14.254	164	592914	20.000	ng	0.00
64) Phenanthrene-d10	17.060	188	1262811	20.000	ng	0.00
76) Chrysene-d12	21.489	240	1556689	20.000	ng	0.00
86) Perylene-d12	24.748	264	1879130	20.000	ng	0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.237	112	1538689	103.448	ng	0.00
7) Phenol-d6	6.813	99	2029858	103.144	ng	0.00
23) Nitrobenzene-d5	8.760	82	1208267	60.437	ng	0.00
42) 2,4,6-Tribromophenol	15.783	330	951096	116.025	ng	0.00
45) 2-Fluorobiphenyl	12.860	172	2409377	54.742	ng	0.00
79) Terphenyl-d14	19.783	244	4624826	53.247	ng	-0.01

Target Compounds	Qvalue

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

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