

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100725\
 Data File : BP025945.D
 Acq On : 07 Oct 2025 16:37
 Operator : RC/JU
 Sample : Q3289-06
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 72-11966

Quant Time: Oct 07 17:01:33 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091225.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 06 19:16:54 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.737	152	150515	20.000	ng	-0.02
21) Naphthalene-d8	10.502	136	548683	20.000	ng	-0.02
39) Acenaphthene-d10	14.354	164	340321	20.000	ng	-0.02
64) Phenanthrene-d10	17.166	188	705313	20.000	ng	-0.02
76) Chrysene-d12	21.607	240	773887	20.000	ng	-0.04
86) Perylene-d12	24.960	264	886824	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.378	112	1178880	126.378	ng	0.00
7) Phenol-d6	6.943	99	1349234	108.819	ng	0.00
23) Nitrobenzene-d5	8.884	82	995764	80.054	ng	-0.02
42) 2,4,6-Tribromophenol	15.878	330	625706	147.404	ng	-0.03
45) 2-Fluorobiphenyl	12.960	172	2175825	85.129	ng	-0.03
79) Terphenyl-d14	19.878	244	3646689	84.771	ng	-0.04

Target Compounds	Qvalue

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

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