

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102225\
 Data File : BP026010.D
 Acq On : 22 Oct 2025 22:42
 Operator : RC/JU
 Sample : Q3402-06
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 41-RA

Quant Time: Oct 23 10:42:13 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP101525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 15 23:48:51 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.696	152	153779	20.000	ng	0.00
21) Naphthalene-d8	10.466	136	619977	20.000	ng	0.00
39) Acenaphthene-d10	14.319	164	459976	20.000	ng	0.00
64) Phenanthrene-d10	17.130	188	1018978	20.000	ng	0.01
76) Chrysene-d12	21.577	240	874935	20.000	ng	0.00
86) Perylene-d12	24.936	264	611463	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.343	112	1330431	141.845	ng	0.00
7) Phenol-d6	6.913	99	1628137	124.386	ng	0.00
23) Nitrobenzene-d5	8.849	82	1211649	87.844	ng	0.00
42) 2,4,6-Tribromophenol	15.842	330	850741	137.386	ng	0.00
45) 2-Fluorobiphenyl	12.925	172	2840144	79.141	ng	0.00
79) Terphenyl-d14	19.854	244	4879285	89.126	ng	0.00

Target Compounds	Qvalue

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

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