

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102225\
 Data File : BP026008.D
 Acq On : 22 Oct 2025 21:20
 Operator : RC/JU
 Sample : Q3399-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COMP-ROLLOFF

Quant Time: Oct 23 10:41:20 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	7.695	152	161883	20.000	ng	0.00
21) Naphthalene-d8	10.460	136	654901	20.000	ng	0.00
39) Acenaphthene-d10	14.325	164	467312	20.000	ng	0.00
64) Phenanthrene-d10	17.136	188	948799	20.000	ng	0.02
76) Chrysene-d12	21.589	240	663420	20.000	ng	0.00
86) Perylene-d12	24.936	264	691050	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.337	112	1184890	120.004	ng	0.00
7) Phenol-d6	6.907	99	1390233	100.894	ng	0.00
23) Nitrobenzene-d5	8.842	82	1135778	77.952	ng	0.00
42) 2,4,6-Tribromophenol	15.860	330	730049	116.045	ng	0.02
45) 2-Fluorobiphenyl	12.931	172	2585253	70.908	ng	0.00
79) Terphenyl-d14	19.854	244	3633726	87.537	ng	0.00

Target Compounds	Qvalue

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

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