

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121225\
 Data File : BP026316.D
 Acq On : 12 Dec 2025 19:07
 Operator : RC/JU
 Sample : Q3839-09
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AN5

Quant Time: Dec 15 02:19:12 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111825.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 15 01:38:31 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.649	152	218390	20.000	ng	-0.02
21) Naphthalene-d8	10.413	136	934074	20.000	ng	-0.03
39) Acenaphthene-d10	14.266	164	637315	20.000	ng	-0.05
64) Phenanthrene-d10	17.078	188	1426387	20.000	ng	-0.05
76) Chrysene-d12	21.536	240	1714188	20.000	ng	-0.04
86) Perylene-d12	24.830	264	1886943	20.000	ng	-0.07
System Monitoring Compounds						
5) 2-Fluorophenol	5.278	112	973904	71.694	ng	-0.01
7) Phenol-d6	6.837	99	1215770	70.903	ng	-0.02
23) Nitrobenzene-d5	8.784	82	733538	44.700	ng	-0.03
42) 2,4,6-Tribromophenol	15.790	330	709812	85.643	ng	-0.04
45) 2-Fluorobiphenyl	12.884	172	1932325	44.081	ng	-0.04
79) Terphenyl-d14	19.825	244	3144226	37.490	ng	-0.05
Target Compounds						
20) 3+4-Methylphenols	8.408	107	44246	2.560	ng	Qvalue 93

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

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