

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111125\
 Data File : BP026105.D
 Acq On : 12 Nov 2025 01:06
 Operator : RC/JU
 Sample : Q3581-08
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RBR200059

Quant Time: Nov 12 01:50:39 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP102925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 30 11:51:34 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.678	152	310420	20.000	ng	-0.01
21) Naphthalene-d8	10.443	136	1185129	20.000	ng	-0.01
39) Acenaphthene-d10	14.313	164	727864	20.000	ng	0.00
64) Phenanthrene-d10	17.137	188	1503498	20.000	ng	0.01
76) Chrysene-d12	21.572	240	1680743	20.000	ng	0.00
86) Perylene-d12	24.895	264	1831494	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.296	112	2483728	132.027	ng	0.00
7) Phenol-d6	6.855	99	2813401	117.498	ng	0.00
23) Nitrobenzene-d5	8.813	82	1929216	101.480	ng	-0.01
42) 2,4,6-Tribromophenol	15.837	330	1369053	173.982	ng	0.00
45) 2-Fluorobiphenyl	12.925	172	4637080	91.544	ng	0.00
79) Terphenyl-d14	19.872	244	7538816	91.129	ng	0.00
Target Compounds						
32) Benzoic acid	9.678	122	7697	4.076	ng	Qvalue 95
84) Bis(2-ethylhexyl)phtha...	21.519	149	9570	2.731	ng	# 97

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

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