

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102225\
 Data File : BP026016.D
 Acq On : 23 Oct 2025 02:46
 Operator : RC/JU
 Sample : Q3368-07
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PRESS-SLUDGE

Quant Time: Oct 23 10:47:17 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	175466	20.000	ng	0.00
21) Naphthalene-d8	10.460	136	751992	20.000	ng	0.00
39) Acenaphthene-d10	14.313	164	463470	20.000	ng	0.00
64) Phenanthrene-d10	17.125	188	720778	20.000	ng	0.00
76) Chrysene-d12	21.577	240	672502	20.000	ng	0.00
86) Perylene-d12	24.918	264	811405	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.337	112	1379928	128.939	ng	0.00
7) Phenol-d6	6.908	99	1780203	119.194	ng	0.00
23) Nitrobenzene-d5	8.843	82	1266001	75.671	ng	0.00
42) 2,4,6-Tribromophenol	15.837	330	588898	94.384	ng	0.00
45) 2-Fluorobiphenyl	12.919	172	2656377	73.462	ng	-0.01
79) Terphenyl-d14	19.842	244	2439194	57.967	ng	-0.01
Target Compounds						
						Qvalue
15) Benzyl Alcohol	7.972	79	40336	3.173	ng	90
20) 3+4-Methylphenols	8.513	107	166114	11.030	ng	93
32) Benzoic acid	9.902	122	24344	2.751	ng	98
84) Bis(2-ethylhexyl)phtha...	21.472	149	64887	2.381	ng	98

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

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