

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040325\
 Data File : BG064171.D
 Acq On : 3 Apr 2025 17:56
 Operator : RC/JU
 Sample : PB167393BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB167393BL

Quant Time: Apr 04 01:44:57 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG030525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 05 15:39:19 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.862	152	25692	20.000	ng	0.00
21) Naphthalene-d8	10.646	136	118561	20.000	ng	0.00
39) Acenaphthene-d10	14.483	164	97336	20.000	ng	0.00
64) Phenanthrene-d10	17.221	188	234504	20.000	ng	0.00
76) Chrysene-d12	21.451	240	246421	20.000	ng	0.00
86) Perylene-d12	24.466	264	258720	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.447	112	226578	137.704	ng	0.00
7) Phenol-d6	7.027	99	311166	139.014	ng	0.00
23) Nitrobenzene-d5	9.007	82	211883	98.760	ng	0.00
42) 2,4,6-Tribromophenol	15.970	330	177238	163.812	ng	0.00
45) 2-Fluorobiphenyl	13.108	172	549402	85.675	ng	0.00
79) Terphenyl-d14	19.847	244	1118423	91.772	ng	0.00

Target Compounds	Qvalue

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

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