

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010422\
 Data File : BP008664.D
 Acq On : 04 Jan 2022 16:25
 Operator : CG/JU
 Sample : PB141778BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB141778BL

Quant Time: Jan 05 04:25:52 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP122821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 30 08:44:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	454982	20.000	ng	-0.01
21) Naphthalene-d8	10.687	136	1785871	20.000	ng	-0.01
39) Acenaphthene-d10	14.516	164	1252093	20.000	ng	-0.01
64) Phenanthrene-d10	17.263	188	3006066	20.000	ng	-0.01
76) Chrysene-d12	21.345	240	3622401	20.000	ng	-0.02
86) Perylene-d12	23.768	264	3967961	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	3613953	144.982	ng	0.00
7) Phenol-d6	7.057	99	4448602	128.217	ng	0.00
23) Nitrobenzene-d5	9.040	82	2623569	87.393	ng	-0.02
42) 2,4,6-Tribromophenol	16.004	330	2679562	158.177	ng	-0.01
45) 2-Fluorobiphenyl	13.145	172	7466576	86.397	ng	-0.01
79) Terphenyl-d14	19.821	244	13332506	75.149	ng	0.00

Target Compounds	Qvalue

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

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