

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031122\
 Data File : BP009360.D
 Acq On : 11 Mar 2022 12:10
 Operator : CG/JU
 Sample : PB143234BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB143234BL

Quant Time: Mar 12 04:49:00 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 05 01:55:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	346484	20.000	ng	0.00
21) Naphthalene-d8	10.616	136	1337448	20.000	ng	0.00
39) Acenaphthene-d10	14.463	164	835467	20.000	ng	0.00
64) Phenanthrene-d10	17.221	188	1713146	20.000	ng	0.00
76) Chrysene-d12	21.339	240	1497612	20.000	ng	# 0.01
86) Perylene-d12	23.756	264	1524506	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	2997904	126.297	ng	0.00
7) Phenol-d6	6.998	99	3648089	120.939	ng	0.01
23) Nitrobenzene-d5	8.975	82	2347362	96.669	ng	0.00
42) 2,4,6-Tribromophenol	15.963	330	1716592	170.157	ng	0.00
45) 2-Fluorobiphenyl	13.086	172	5674727	93.408	ng	0.00
79) Terphenyl-d14	19.798	244	7961633	100.324	ng	0.00

Target Compounds	Qvalue

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

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