

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031522\
 Data File : BP009399.D
 Acq On : 15 Mar 2022 15:39
 Operator : CG/JU
 Sample : PB143311BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB143311BL

Quant Time: Mar 16 01:56:42 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	7.811	152	307256	20.000	ng	-0.01
21) Naphthalene-d8	10.605	136	1161883	20.000	ng	-0.01
39) Acenaphthene-d10	14.457	164	727463	20.000	ng	0.00
64) Phenanthrene-d10	17.216	188	1422246	20.000	ng	0.00
76) Chrysene-d12	21.327	240	1243146	20.000	ng	# 0.00
86) Perylene-d12	23.739	264	1370883	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	2568030	121.999	ng	0.01
7) Phenol-d6	7.005	99	3137762	117.302	ng	0.02
23) Nitrobenzene-d5	8.969	82	2038609	96.640	ng	0.00
42) 2,4,6-Tribromophenol	15.951	330	1525355	173.649	ng	0.00
45) 2-Fluorobiphenyl	13.075	172	4996427	94.453	ng	-0.01
79) Terphenyl-d14	19.792	244	6784069	102.984	ng	0.00

Target Compounds	Qvalue

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

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