

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013122\
 Data File : BP008962.D
 Acq On : 31 Jan 2022 13:24
 Operator : CG/JU
 Sample : N1303-08
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CC-8

Quant Time: Feb 01 04:19:02 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 01 04:16:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.828	152	272724	20.000	ng	0.00
21) Naphthalene-d8	10.628	136	1029475	20.000	ng	0.00
39) Acenaphthene-d10	14.469	164	692119	20.000	ng	0.00
64) Phenanthrene-d10	17.228	188	1503518	20.000	ng	0.00
76) Chrysene-d12	21.316	240	1451883	20.000	ng	-0.01
86) Perylene-d12	23.727	264	1290145	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	2161948	122.588	ng	0.00
7) Phenol-d6	7.016	99	2379359	111.414	ng	0.00
23) Nitrobenzene-d5	8.993	82	1811139	99.881	ng	0.00
42) 2,4,6-Tribromophenol	15.969	330	1579899	156.821	ng	0.00
45) 2-Fluorobiphenyl	13.093	172	4684097	89.249	ng	0.00
79) Terphenyl-d14	19.792	244	7617372	88.559	ng	0.00
Target Compounds						
32) Benzoic acid	9.946	122	20255	2.150	ng	Qvalue 96

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

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