

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020222\
 Data File : BP009010.D
 Acq On : 02 Feb 2022 17:11
 Operator : CG/JU
 Sample : N1217-13
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 VNJ-242

Quant Time: Feb 03 06:25:10 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 03 06:20:22 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	187463	20.000	ng	0.00
21) Naphthalene-d8	10.622	136	702355	20.000	ng	0.00
39) Acenaphthene-d10	14.463	164	441012	20.000	ng	0.00
64) Phenanthrene-d10	17.216	188	930384	20.000	ng	-0.01
76) Chrysene-d12	21.316	240	1008952	20.000	ng	0.00
86) Perylene-d12	23.722	264	1128795	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.423	112	1448366	119.478	ng	0.01
7) Phenol-d6	7.011	99	1581814	107.757	ng	0.00
23) Nitrobenzene-d5	8.981	82	1253752	101.345	ng	0.00
42) 2,4,6-Tribromophenol	15.963	330	981648	152.919	ng	0.00
45) 2-Fluorobiphenyl	13.087	172	3118262	93.244	ng	0.00
79) Terphenyl-d14	19.786	244	4915739	82.239	ng	0.00
Target Compounds						
15) Benzyl Alcohol	8.081	79	63531	6.169	ng	Qvalue 94
22) Acetophenone	8.646	105	99700	5.572	ng	# 92
32) Benzoic acid	9.952	122	40588	6.314	ng	96

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

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