

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012423\
 Data File : BP013565.D
 Acq On : 24 Jan 2023 11:43
 Operator : CG/JU
 Sample : 01233-10
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 WC-38

Quant Time: Jan 25 03:46:33 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP012323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 25 03:43:27 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.164	152	253707	20.000	ng	0.00
21) Naphthalene-d8	10.993	136	972664	20.000	ng	# 0.00
39) Acenaphthene-d10	14.793	164	544889	20.000	ng	0.00
64) Phenanthrene-d10	17.545	188	1077804	20.000	ng	0.00
76) Chrysene-d12	21.622	240	971180	20.000	ng	0.00
86) Perylene-d12	24.210	264	1162665	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.687	112	1726783	116.615	ng	0.00
7) Phenol-d6	7.299	99	2092403	104.992	ng	0.00
23) Nitrobenzene-d5	9.334	82	1256986	85.314	ng	0.00
42) 2,4,6-Tribromophenol	16.281	330	714060	130.649	ng	0.00
45) 2-Fluorobiphenyl	13.422	172	2650143	66.761	ng	0.00
79) Terphenyl-d14	20.081	244	3436443	63.656	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	11.046	128	2262982	43.276	ng	97
37) 2-Methylnaphthalene	12.634	142	1300841	34.998	ng	99
38) 1-Methylnaphthalene	12.851	142	470768	13.479	ng	99
71) Phenanthrene	17.587	178	145625	2.428	ng	98

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

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