

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013023\
 Data File : BP013659.D
 Acq On : 30 Jan 2023 23:10
 Operator : CG/JU
 Sample : 01329-12
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COMP-8-BORING-2-21-28

Quant Time: Jan 31 04:45:22 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP012623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 31 04:38:43 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.146	152	223368	20.000	ng	0.00
21) Naphthalene-d8	10.975	136	787371	20.000	ng	0.00
39) Acenaphthene-d10	14.781	164	455839	20.000	ng	0.00
64) Phenanthrene-d10	17.533	188	975415	20.000	ng	0.00
76) Chrysene-d12	21.610	240	948730	20.000	ng	-0.01
86) Perylene-d12	24.186	264	1197792	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.681	112	1763897	141.508	ng	0.00
7) Phenol-d6	7.293	99	2041671	121.249	ng	0.00
23) Nitrobenzene-d5	9.316	82	1294540	90.341	ng	-0.01
42) 2,4,6-Tribromophenol	16.269	330	925283	133.324	ng	0.00
45) 2-Fluorobiphenyl	13.404	172	3329795	104.328	ng	0.00
79) Terphenyl-d14	20.063	244	5320043	98.953	ng	0.00
Target Compounds						
						Qvalue
74) Di-n-butylphthalate	18.463	149	7697	3.741	ng	# 94
84) Bis(2-ethylhexyl)phtha...	21.492	149	14877	4.384	ng	97
85) Di-n-octyl phthalate	22.474	149	41660	3.198	ng	96
88) Benzo(b)fluoranthene	23.392	252	169197	2.200	ng	# 97
89) Benzo(k)fluoranthene	23.445	252	174766	2.289	ng	# 96
92) Benzo(g,h,i)perylene	27.739	276	144878	2.100	ng	# 91

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

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